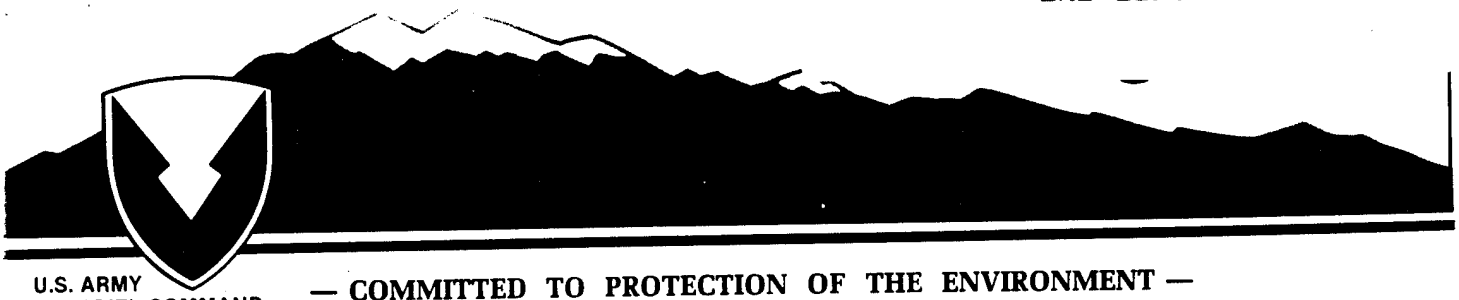

U.S. ARMY
MATERIEL COMMAND

— COMMITTED TO PROTECTION OF THE ENVIRONMENT —

FINAL PHASE II
DATA ADDENDUM
SITE 36-7: SOLID WASTE BURIAL/SANITARY PI

September 1988
Contract Number DAAK11-84-D-0016
(Version 3.1)

Rocky

Environmental Science And Engineering, Inc.
Harding Lawson Associates

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THE PHASE II PROGRAM FOR SITE 36-7 CONSISTED OF 8 PIT BORINGS YIELDING 22 SAMPLES AND 22 SOIL BORINGS YIELDING 40 SAMPLES. THE PIT BORING LOCATIONS WERE SELECTED TO EXPLORE POTENTIAL DISPOSAL TRENCHES. SELECTED SAMPLES WERE ANALYZED FOR VOLATILE AND SEMIVOLATILE ORGANICS, ARMY AGENT DEGRADATION PRODUCTS, AND METALS. DLDRN, ENDRN, ISODR, CPMS, CPMSO, CPMSO2, PB, ZN, AS, HG, FC2A, AND SEVERAL TENTATIVELY IDENTIFIED NONTARGET COMPOUNDS WERE DETECTED.

RESULTS OF THE PHASE II SAMPLING PROGRAM WILL BE ASSESSED AS PART OF THE OVERALL ANALYSIS FOR THE CENTRAL STUDY AREA REPORT.

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Rocky Mountain Arsenal

FINAL PHASE II
DATA ADDENDUM
SITE 36-7: SOLID WASTE BURIAL/SANITARY PIT

September 1988
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(Version 3.1)

PREPARED BY

ENVIRONMENTAL SCIENCE AND ENGINEERING, INC.
Harding Lawson Associates

PREPARED FOR

U.S. ARMY PROGRAM MANAGER'S OFFICE FOR THE ROCKY MOUNTAIN ARSENAL

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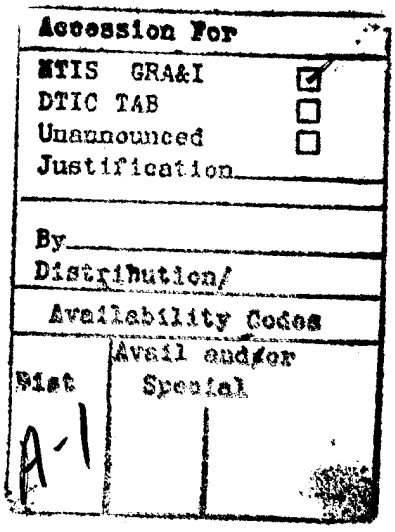

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SITE 36-7: SOLID WASTE BURIAL/SANITARY PIT

1.0 PHASE_II_PROGRAM

As a result of the Phase I Contamination Assessment at Rocky Mountain Arsenal (RMA), a Phase II program was initiated at Site 36-7 in February, 1988. The Phase II Program was generally conducted as presented in the Phase I Contamination Assessment Report (CAR) (ESE, 1988, RIC#88067R07), except that only six samples of trench material were obtained instead of the eight samples planned.

The Phase II investigation of Site 36-7 consisted of 8 pit borings yielding 22 samples and 22 soil borings yielding 40 samples. Pit boring locations were selected to explore potential disposal trenches, and soil boring sites were selected to investigate the remaining areas within Site 36-7. Observation pits were excavated with a backhoe in the eight suspected disposal trenches as defined by geophysical anomalies, interpretation of aerial photography, and review of historical data. Each pit was excavated to the trench bottom, as determined by visual inspection. A grab sample believed to be representative of the most contaminated material within the trench was obtained from excavated material. The pit was backfilled with clean (metal free) material, and the location was staked. The pit borings were drilled 5 feet (ft) below the base of the suspected disposal trench and sampled at the 0- to 1- and 4- to 5-ft intervals below the base.

A continuous shallow trench was excavated with a backhoe in the southernmost anomaly to identify the locations and orientations of suspected disposal trenches and to allow observation pits and pit borings to be targeted more accurately.

All soil samples (except for grab samples) were collected using the continuous soil sampling method detailed in the Task 1 Technical Plan (ESE, 1985, RIC#85127R07). Samples were obtained at predetermined intervals unless field conditions (e.g., depth to water table, staining, etc.) necessitated adjustment in the interval selections. Seven 1-ft-deep borings were drilled and sampled using hand-auger equipment. Locations of the observation pits and pit borings are shown on Figure 36-7-II-1 (See Section 2.0),

and locations of all Phase II borings are shown on Figure 36-7-II-2 (See Section 4.0).

The actual Phase II investigation varied from the proposed Phase II program in that only six grab samples of disposal trench material were obtained instead of the eight samples planned. A grab sample from Pit Boring 3682 was not obtained, because only natural soil and bedrock were encountered at that location. A grab sample from Pit Boring 3687 was not obtained, because the excavation caved and resulted in unsafe conditions for equipment and personnel. Boring 3687, however, was drilled through the trench material. The depth of the trench was determined from Boring 3687, and soil samples were obtained from beneath the trench.

Prior to any Phase II drilling, the Program Manager's Office (PMO), Environmental Science and Engineering (ESE), Morrison-Knudsen Engineers (MKE), and Harding Lawson Associates (HLA) formulated procedures for MKE to obtain subsamples from selected soil cores during Phase II drilling. MKE did not request subsamples at Site 36-7.

The following table summarizes the Phase II investigation at Site 36-7:

Boring No.	Total Depth (ft)	Sampling Interval(s)	No. of Samples
3683*	10	5-6, 9-10	2
3684*	15	2-3+, 10-11, 14-15	3
3685*	17	7-8+, 10-11, 15-16	3
3686*	15	7-8+, 10-11, 14-15	3
3687*	22	17-18, 21-22	2
3688*	14	5-6+, 9-10, 13-14	3
3689*	16	3-4+, 11-12, 15-16	3
3690*	15	4-5+, 10-11, 14-15	3
3691	5	0-1, 4-5	2
3692	5	0-1, 4-5	2
3693	5	0-1, 4-5	2
3694	5	0-1, 4-5	2
3695	5	0-1, 4-5	2
3696	5	0-1, 4-5	2
3697	5	0-1, 4-5	2
3698	5	0-1, 4-5	2
3699	5	0-1, 4-5	2
3700	1	0-1	1
3701	1	0-1	1

3702	1	0-1	1
3703	1	0-1	1
3704	1	0-1	1
3705	1	0-1	1
3706	1	0-1	1
3707	10	0-1, 4-5, 9-10	3
3708	10	0-1, 4-5, 9-10	3
3709	10	0-1, 4-5, 9-10	3
3710	3	0-1, 2-3	2
3711	3	0-1, 2-3	2
3712	3	0-1, 2-3	2
		Total	62

* Pit boring
+ Grab sample of trench material

The Phase II analytical program was conducted as planned with the exception of the grab samples from Borings 3683 and 3687, which were not collected. Fifty-six samples were analyzed for semivolatile organic (SVO) compounds by gas chromatography/mass spectrometry (GC/MS) and for arsenic and mercury by atomic absorption (AA). Sixty-two samples were analyzed for cadmium, chromium, copper, lead, and zinc by the inductively coupled argon plasma (ICP) method. The 4-to 5-ft samples below the base of disposal trenches from Borings 3683 to 3690 and the 9- to 10-ft samples from Borings 3707 through 3709 were analyzed for volatile organic (VO) compounds by GC/MS. Six samples were analyzed by high-performance liquid chromatography (HPLC) for the Army Agent Degradation Products (ADP) thiodiglycol (TDGCL) and chloroacetic acid (CLC2A), and by ion chromatography (IONCHROM) for fluoro-acetic acid (FC2A), isopropylmethyphosphonic acid (IMPA), and methylphosphonic acid (MPA).

The six samples from the Phase II borings triangulated around Phase I Boring 3111 were analyzed for ICP metals to investigate the elevated cadmium in Boring 3111. The remaining samples were analyzed for the Phase I suite of analytes, because Phase I samples were not collected in the geophysical anomalies and disposal trenches investigated in the Phase II program. Selected Phase II samples were also analyzed for IMPA and TDGCL to screen for ADPs, as the HPLC and IONCHROM methods were not available during the Phase II program.

Phase I and Phase II analytical methods for Site 36-7 samples were the same for VO and SVO compounds, ICP metals, arsenic, and mercury; therefore, Phase I and Phase II results are directly comparable. Appendix 36-7-II-A provides a complete list of analytes, analytical methods, and standard abbreviations used in the Phase I and Phase II investigations.

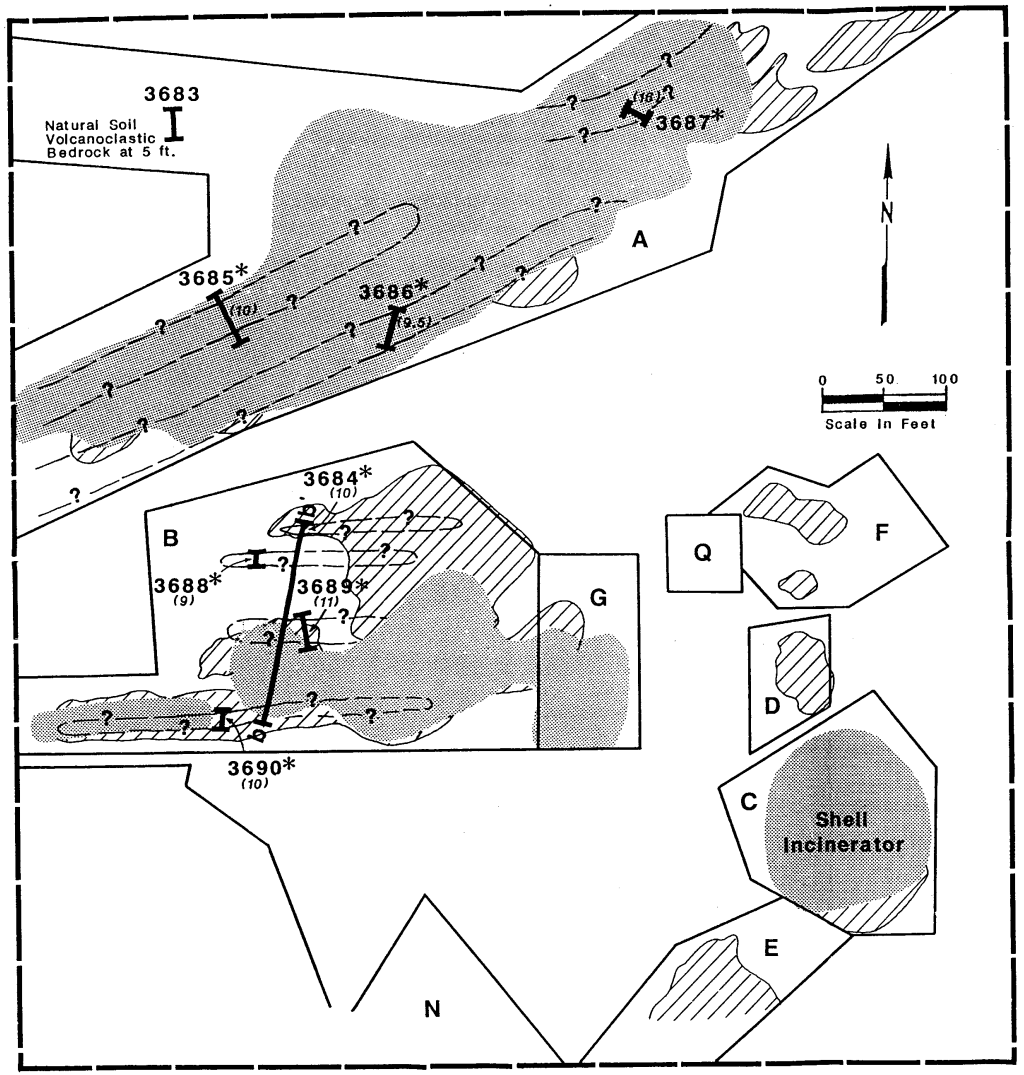
2.0 PHASE II FIELD OBSERVATIONS

Surface conditions at Site 36-7 have not changed appreciably since the Phase I investigation was completed in the summer of 1985 (ESE, 1988, RIC#88063R07). Figure 36-7-II-1 shows the locations of excavated pits and trenches, excavation depths, trend outlines of possible disposal trenches as interpreted from aerial photographs, actual widths of intercepted disposal trenches, and a cross section showing the continuous shallow excavation profile with the intercepted disposal trenches projected onto the cross section. Table 36-7-II-1 presents a summary of pit excavation activities and observations.

Ground water was not encountered at any of the Phase II pits or borings at Site 36-7. Volcaniclastic bedrock was encountered in seven of the eight pit borings. Bedrock depths are presented on Table 36-7-II-1.

For safety purposes, air monitoring was conducted using a photoionization detector (PID) during drilling and excavation activities. In general, PID readings in the auger annulus and open trenches were background except in Pit Boring 3687, which registered a PID reading of 13.2 in the hollow-stem annulus at the 16- to 18-ft depth, and Boring 3698, which had a PID reading of 24 in the 0- to 1-ft sample. PID readings were at background levels in the breathing zone during field activities.

An M18A2 test kit was used to detect the presence of chemical agents in trenches, boreholes, and soil samples. Specifically at RMA, the M18A2 test kit is used to detect Sarin (GB), nerve agent (VX), mustard (H), and Lewisite (L), based on the knowledge that these agents were manufactured, stored, or demilitarized at the site. The detection limit for mustard agents is 0.5 milligrams per cubic meter (mg/m^3), and the detection limit for GB, VX, and L is 0.2 mg/m^3. The detection limits for L and VX in soil are 5 and 5.9 parts per million (ppm), respectively. All M18A2 field test results for chemical agents at this site were negative.



LEGEND

(9) Excavation Location
Depth to Bottom of Disposal Trench Material
Intercepted Buried Trench

—?— Possible Trench Boundary Inferred from Aerial Photography

3688* Refer to Table 36-7-II-1, for Description of Disposal Trench Contents

——— Geophysical Anomaly; **A, B, C,** etc.

GEOPHYSICAL KEY

Magnetic Anomaly

Combined Electromagnetic and Magnetic Anomaly

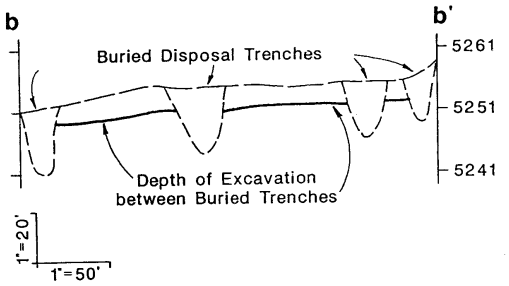


Figure 36-7-II-1

SITE 36-7, TRENCH EXLORATION INVESTIGATION
PHASE II RESULTS

SOURCE: HLA, 1988

Prepared for:
U.S. Army Program Manager's Office
For Rocky Mountain Arsenal

Aberdeen Proving Ground, Maryland

Table 36-7-II-1. Summary of Pit Boring Activities and Observations

Pit Boring No.	Disposal Trench Depth (ft)	Pit Boring Depth (ft)	Disposal Trench Contents	Comments
3683	Trench not encountered	10	None	Volcaniclastic bedrock at 5 ft
3684	10	15	Paper, plastic, metal, lumber	Siltstone/ volcaniclastic bedrock at 12.5 ft
3685	10	17	Plastic pipe, scrap metal, wood	Volcaniclastic bedrock at 9.5 ft
3686	9.5	15	Paper, cans, respirators, M-19 casings, scrap metal	Volcaniclastic bedrock at 9.2 ft
3687	16*	22	Paper, lumber, plastic	Excavation pit abandoned due to severe instability
3688	9	14	Paper, lumber, plastic	Volcaniclastic bedrock at 10.4 ft
3689	11	16	Old tires, auto-motive trash	Volcaniclastic bedrock at 15.5 ft
3690	10	15	Paper, wood, plastic	Volcaniclastic bedrock at 10 ft

* Estimated from soil core obtained during drilling

Source: ESE, 1988.

Samples at Site 36-7 were also tested for chemical agents by the RMA laboratory, because historical evidence indicated possible agent presence. A composite of aliquots from each sample was initially analyzed for GB, VX, H, and L. If agent had been detected, individual samples from each boring would have been analyzed to identify stratigraphic location. No positive results for chemical agents were found at this site.

3.0 PHASE II GEOPHYSICAL EXPLORATION

No geophysical survey was conducted at Site 36-7 for the Phase II investigation. Fifteen boring locations were cleared, however, for safety purposes in accordance with the borehole clearance geophysical program detailed in the Task 1 Technical Plan (ESE, 1985, RIC#85127R07). Borehole site clearance was used to ensure that drilling would not encounter buried unexploded ordnance or other metal that could pose a significant safety risk. Magnetic intensity readings were obtained with a gradiometer. A 20-ft-square grid was centered at each boring location, and gradiometer readings were obtained at 5-ft intervals throughout the area. A contour map was prepared from the data and was used to place the boring in the safest location within the geophysical plot. Following borehole site clearance, a metal detector was used to check for surficial (0 to 2 ft) metal. None of the 15 borings were relocated as a result of the borehole site clearance conducted at Site 36-7.

The seven hand-augered boring locations were cleared by a metal detector survey for shallow (0 to 2 ft) buried metal. None of these borings had to be relocated after the metal detector survey.

4.0 PHASE II ANALYTE LEVELS AND DISTRIBUTION

Table 36-7-II-2 contains indicator ranges and a statistical summary of Phase II analytical results. A summary of analytical data for each sample, including lithology and air monitoring results, is presented in Table 36-7-II-3. A tabulation of all analytical data associated with the Phase II investigation at this site is presented in Appendix 36-7-II-B.

To assess the significance of metal and organic analytical values, indicator ranges were established during the Phase I program. For organic compounds, the indicator level is the method detection limit. For metals, a range of values was chosen to reflect the upper end of the expected natural range for each metal as normally found in RMA alluvial soil. The procedure for establishing indicator ranges is presented in the Introduction to the Contamination Assessment Reports (ESE, 1987, RIC#88204R02). Concentrations within or above indicator ranges for Phase I and Phase II data are presented in Figure 36-7-II-2.

Eleven samples were analyzed for VO compounds; the only target compound detected was methylene chloride in two samples at low concentrations (Boring 3684, 14 to 15 ft, 1 ppm; Boring 3690, 14 to 15 ft, 0.7 ppm). Fifty-six samples were analyzed for SVO; dieldrin was the most prevalent compound detected with concentrations ranging from 0.3 to 10 parts per million (ppm) in 8 samples. Four samples contained chlorophenylmethyl sulfone ($CPMSO_2$) at concentrations ranging from 0.7 to 2 ppm. Isodrin, endrin, chlorophenylmethyl sulfide (CPMS), and chlorophenylmethyl sulfoxide (CPMSO) were each detected in one sample.

Four of the dieldrin detections, the endrin detection (1 ppm), and the isodrin detection (0.5 ppm) were found in samples from the pit borings. Phase II results showed a dieldrin detection of 10 ppm in the grab sample from the 2-to 3-ft interval of Pit Boring 3684. Lower levels of dieldrin were detected in Pit Borings 3686 (1 ppm, 7- to 8-ft interval), 3688 (0.4 ppm, 5- to 6-ft interval), and 3691 (2 ppm, 0- to 1-ft interval). The highest concentration was detected in Pit Boring 3684, which is approximately 50 ft northwest of Boring 3113. As noted in Table 36-7-II-1, Pit Boring 3684 (2- to 3-ft interval) was sampled from a buried disposal trench containing paper, lumber, metal, and plastic.

Table 36-7-II-2. Summary of Analytical Results for Site 36-7 Phase II Soil Samples

Constituent	Number of Samples*	CONCENTRATIONS (ug/g) Range	Mean**	Median**	Standard Deviation**	ESE Detection Limit	Indicator Level
IMPA (N=6)+							
IMPA	0	--	--	--	--	2.1	DL
Fluoroacetic Acid	2	2.9-15	--	--	--	2.0	DL
MPA	0	--	--	--	--	2.0	DL
TDGCL (N=6)+							
Thiodiglycol	0	--	--	--	--	2.6	DL
Chloroacetic Acid	0	--	--	--	--	18	DL
VOLATILE ORGANICS (N=11)+							
Methylene Chloride	2	0.7-1	--	--	--	0.3	DL
SEMIVOLATILE ORGANICS (N=56)+							
CMPS	1	8	--	--	--	0.3	DL
CMPSO	1	4	--	--	--	0.4	DL
CMPSO2	4	0.7-2	--	--	--	0.3	DL
Dieldrin	8	0.3-10	2	0.9	3	0.3	DL
Endrin	1	1	--	--	--	0.7	DL
Isodrin	1	0.5	--	--	--	0.3	DL
ICP METALS (N=62)+							
Cadmium	1	1.3	--	--	--	0.92	DL-2.0
Chromium	37	8.4-18	12	12	2.5	7.2	25-40
Copper	62	6.5-77	16	11	13	4.8	20-35
Lead	7	22-81	50	48	23	17	25-40
Zinc	62	25-250	56	44	37	16	60-80
ARSENIC (N=56)+	7	6.0-15	7.9	6.8	3.3	4.7	DL-10
MERCURY (N=56)+	14	0.054-2.1	0.38	0.11	0.64	0.050	DL-0.10

* Number of samples in which constituent was detected. Only these sample results were used in statistical analyses.
** Statistics not calculated when constituent detected in fewer than five samples.
\+ Number of samples analyzed by laboratory.
DL Detection limit.

Source: ESE, 1988.

Table 36-7-II-3. Concentrations of Target Analtyes Above Detection Limits in Site 36-7 Phase II Soil Samples (page 1 of 5)

Boring Number	3683	3683	3684	3684	3684	3685	3685	3685	3686	3686	3686	3687
Depth (ft)	5-6	9-10	2-3	10-11	14-15	7-8	10-11	15-16	7-8	10-11	14-15	17-18
Geologic Material	Volcani-clastic (Denver Fm)	Volcani-clastic (Denver Fm)	Trench Material	Clayey Silt	Volcani-clastic (Denver Fm)	Trench Material	Volcani-clastic (Denver Fm)	Volcani-clastic (Denver Fm)	Trench Material	Volcani-clastic (Denver Fm)	Volcani-clastic (Denver Fm)	Clayey Silt
AIR MONITORING												
PID*	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	13.2
SOIL CHEMISTRY												
IMPA (ug/g)												
IMPA	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ
Fluoroacetic Acid	NRQ	NRQ	BDL	NRQ	NRQ	15	NRQ	NRQ	BDL	NRQ	NRQ	NRQ
MPA	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ
TDGCL (ug/g)												
Thiodiglycol	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ
Chloroacetic Acid	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ
Volatile Organics (VO) by GC/MS (ug/g)												
Methylene Chloride	NRQ	BDL	NRQ	NRQ	1	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ
Semivolatile Organics (SVO) by GC/MS (ug/g)												
CPMS	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
CPMSO	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
CPMSO2	BDL	BDL	BDL	BDL	BDL	0.7	BDL	BDL	BDL	BDL	1	BDL
Dieldrin	BDL	BDL	10	0.3	BDL	BDL	BDL	BDL	1	BDL	BDL	BDL
Endrin	BDL	BDL	1	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Isodrin	BDL	BDL	0.5	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
ICP metals (ug/g)												
Cadmium	BDL	BDL	BDL	BDL	BDL	1.3	BDL	BDL	BDL	BDL	BDL	BDL
Chromium	BDL	BDL	17	9.5	BDL	12	BDL	BDL	12	BDL	BDL	BDL
Copper	34	32	25	23	51	33	35	41	18	12	20	27
Lead	BDL	BDL	42	48	BDL	61	BDL	BDL	BDL	BDL	BDL	BDL
Zinc	67	71	76	49	87	170	65	85	62	70	55	58
Arsenic (ug/g)	BDL	BDL	7.5	BDL	BDL	7.0	BDL	BDL	BDL	BDL	BDL	BDL
Mercury (ug/g)	BDL	BDL	0.11	0.14	BDL	1.7	0.075	BDL	BDL	BDL	BDL	BDL

< Higher detection limit due to dilution or soil matrix masking effects.
> Quantitative concentration was not achieved due to dilution constraints.
* As calibrated to an isobutylene standard.
BDL Below detection limit.
BKD No reading above ambient background.
NRQ Analysis not requested.
NA Not analyzed.

Table 36-7-II-3. Concentrations of Target Analtyes Above Detection Limits in Site 36-7 Phase II Soil Samples (page 2 of 5)

Boring Number	3687	3688	3688	3688	3689	3689	3689	3690	3690	3690	3691	3691
Depth (ft)	21-22	5-6	9-10	13-14	3-4	11-12	15-16	4-5	10-11	14-15	0-1	4-5
Geologic Material	Silty Clay	Trench Material	Clayey Silt	Volcani-clastic (Denver Fm)	Trench Material	Sandy Silt	Silt	Trench Material	Volcani-clastic (Denver Fm)	Volcani-clastic (Denver Fm)	Silty Sand	Silty Sand
AIR MONITORING												
PID*	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD
SOIL CHEMISTRY												
IMPA (ug/g)												
IMPA	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ	NRQ
Fluoroacetic Acid	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	2.9	NRQ	NRQ	NRQ	NRQ
MPA	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ	NRQ
TDGCL (ug/g)												
Thiodiglycol	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ	NRQ
Chloroacetic Acid	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ	NRQ
Volatile Organics (VO) by GC/MS (ug/g)												
Methylene Chloride	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	0.7	NRQ	NRQ
Semivolatile Organics (SVO) by GC/MS (ug/g)												
CPMS	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	8
CPMSO	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	4
CPMSO2	BDL	BDL	BDL	BDL	BDL	2	2	BDL	BDL	BDL	BDL	BDL
Dieldrin	BDL	0.4	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	2	BDL
Endrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Isodrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
ICP metals (ug/g)												
Cadmium	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Chromium	BDL	15	BDL	BDL	BDL	14	9.9	17	9.4	12	12	12
Copper	34	77	11	41	21	13	17	20	10	15	11	9.3
Lead	BDL	BDL	BDL	BDL	81	BDL	BDL	22	BDL	BDL	BDL	BDL
Zinc	85	100	39	78	250	48	57	150	40	57	44	41
Arsenic (ug/g)	BDL	6.8	BDL	BDL	BDL	BDL	BDL	6.7	BDL	BDL	BDL	BDL
Mercury (ug/g)	0.063	0.30	0.13	BDL	0.11	BDL	BDL	BDL	BDL	BDL	BDL	BDL

< Higher detection limit due to dilution or soil matrix masking effects.
> Quantitative concentration was not achieved due to dilution constraints.
* As calibrated to an isobutylene standard.
BDL Below detection limit.
BKD No reading above ambient background.
NRQ Analysis not requested.
NA Not analyzed.

Table 36-7-II-3. Concentrations of Target Analtyes Above Detection Limits in Site 36-7 Phase II Soil Samples (page 3 of 5)

Boring Number	3692	3692	3693	3693	3694	3694	3695	3695	3696	3696	3697	3697
Depth (ft)	0-1	4-5	0-1	4-5	0-1	4-5	0-1	4-5	0-1	4-5	0-1	4-5
Geologic Material	Silty Sand	Silty Sand	Silty Sand	Silty Sand	Silty Sand	Silty Sand	Sandy Silt	Sandy Silt	Silty Sand	Silty Sand	Silty Sand	Silty Sand
AIR MONITORING												
PID*	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD
SOIL CHEMISTRY												
IMPA (ug/g)												
IMPA	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Fluoroacetic Acid	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
MPA	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
TDGCL (ug/g)												
Thiodiglycol	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Chloroacetic Acid	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Volatile Organics (VO) by GC/MS (ug/g)												
Methylene Chloride	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Semivolatile Organics (SVO) by GC/MS (ug/g)												
CPMS	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
CPMSO	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
CPMSO2	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Dieldrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Endrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Isodrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
ICP metals (ug/g)												
Cadmium	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Chromium	BDL	BDL	BDL	12	11	9.0	18	15	BDL	12	9.3	BDL
Copper	8.3	8.8	8.9	13	14	9.1	13	15	10	8.4	8.8	6.8
Lead	BDL	BDL	BDL	BDL	72	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Zinc	35	35	32	50	130	36	56	56	39	36	45	33
Arsenic (ug/g)	BDL	BDL	BDL	BDL	15	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Mercury (ug/g)	BDL	BDL	BDL	BDL	BDL	BDL	BDL	2.1	BDL	0.061	BDL	BDL

< Higher detection limit due to dilution or soil matrix masking effects.
> Quantitative concentration was not achieved due to dilution constraints.
* As calibrated to an isobutylene standard.
BDL Below detection limit.
BKD No reading above ambient background.
NRQ Analysis not requested.
NA Not analyzed.

Table 36-7-II-3. Concentrations of Target Analtyes Above Detection Limits in Site 36-7 Phase II Soil Samples (page 4 of 5)

Boring Number	3698	3698	3699	3699	3700	3701	3702	3703	3704	3705	3706	3707	3707
Depth (ft)	0-1	4-5	0-1	4-5	0-1	0-1	0-1	0-1	0-1	0-1	0-1	0-1	4-5
Geologic Material	Silty Sand	Silty Sand	Silty Sand	Silty Sand	Sandy Silt	Sandy Silt	Sandy Silt	Sandy Silt	Silty Sand	Sandy Silt	Sandy Silt	Silty Sand	Sand
AIR MONITORING													
PID*	24	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD
SOIL CHEMISTRY													
IMPA (ug/g)													
IMPA	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Fluoroacetic Acid	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
MPA	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
TDGCL (ug/g)													
Thiodiglycol	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Chloroacetic Acid	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Volatile Organics (VO) by GC/MS (ug/g)													
Methylene Chloride	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Semivolatile Organics (SVO) by GC/MS (ug/g)													
CPMS	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
CPMSO	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
CPMSO2	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Dieldrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	4	BDL	BDL	BDL	BDL	BDL
Endrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Isodrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
ICP metals (ug/g)													
Cadmium	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Chromium	BDL	9.5	BDL	11	8.8	BDL	11	12	16	10	12	BDL	11
Copper	7.4	8.9	7.7	10	9.8	8.4	12	13	8.7	11	9.3	7.0	9.4
Lead	BDL	BDL	BDL	BDL	BDL	BDL	23	BDL	BDL	BDL	BDL	BDL	BDL
Zinc	34	39	35	44	37	28	49	46	33	52	39	25	35
Arsenic (ug/g)	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	6.2	BDL	BDL
Mercury (ug/g)	BDL	BDL	BDL	BDL	0.054	0.12	0.32	0.065	BDL	BDL	BDL	BDL	BDL

< Higher detection limit due to dilution or soil matrix masking effects.
> Quantitative concentration was not achieved due to dilution constraints.
* As calibrated to an isobutylene standard.
BDL Below detection limit.
BKD No reading above ambient background.
NRQ Analysis not requested.
NA Not analyzed.

Table 36-7-II-3. Concentrations of Target Analtyes Above Detection Limits in Site 36-7 Phase II Soil Samples (page 5 of 5)

Boring Number	3707	3708	3708	3708	3709	3709	3709	3710	3710	3711	3711	3712	3712
Depth (ft)	9-10	0-1	4-5	9-10	0-1	4-5	9-10	0-1	2-3	0-1	2-3	0-1	2-3
Geologic Material	Silty Sand	Sandy Silt	Clayey Silt	VFG Sand	Sandy Silt	Clayey Silt	VFG Sand	Silty Sand	Silty Sand	Silty Sand	Silty Sand	Silty Sand	Silty Sand
AIR MONITORING													
PID*	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD	BKD
SOIL CHEMISTRY													
IMPA (ug/g)													
IMPA	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Fluoroacetic Acid	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
MPA	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
TDGCL (ug/g)													
Thiodiglycol	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Chloroacetic Acid	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Volatile Organics (VO) by GC/MS (ug/g)													
Methylene Chloride	BDL	NRQ	NRQ	BDL	NRQ	NRQ	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Semivolatile Organics (SVO) by GC/MS (ug/g)													
CPMS	BDL	BDL	BDL	BDL	BDL	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
CPMSO	BDL	BDL	BDL	BDL	BDL	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
CPMSO2	BDL	BDL	BDL	BDL	BDL	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Dieldrin	BDL	0.5	BDL	BDL	0.4	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Endrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Isodrin	BDL	BDL	BDL	BDL	BDL	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
ICP metals (ug/g)													
Cadmium	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Chromium	9.3	BDL	9.1	16	12	15	11	BDL	BDL	11	10	BDL	11
Copper	11	11	9.8	12	14	14	13	8.8	6.5	9.1	7.8	7.5	9.0
Lead	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL	BDL
Zinc	34	39	35	39	54	51	37	34	28	37	32	34	40
Arsenic (ug/g)	BDL	BDL	BDL	BDL	BDL	BDL	6.0	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ
Mercury (ug/g)	BDL	BDL	BDL	BDL	BDL	BDL	BDL	NRQ	NRQ	NRQ	NRQ	NRQ	NRQ

< Higher detection limit due to dilution or soil matrix masking effects.
> Quantitative concentration was not achieved due to dilution constraints.
* As calibrated to an isobutylene standard.
BDL Below detection limit.
BKD No reading above ambient background.
NRQ Analysis not requested.
NA Not analyzed.

Source: ESE, 1988.

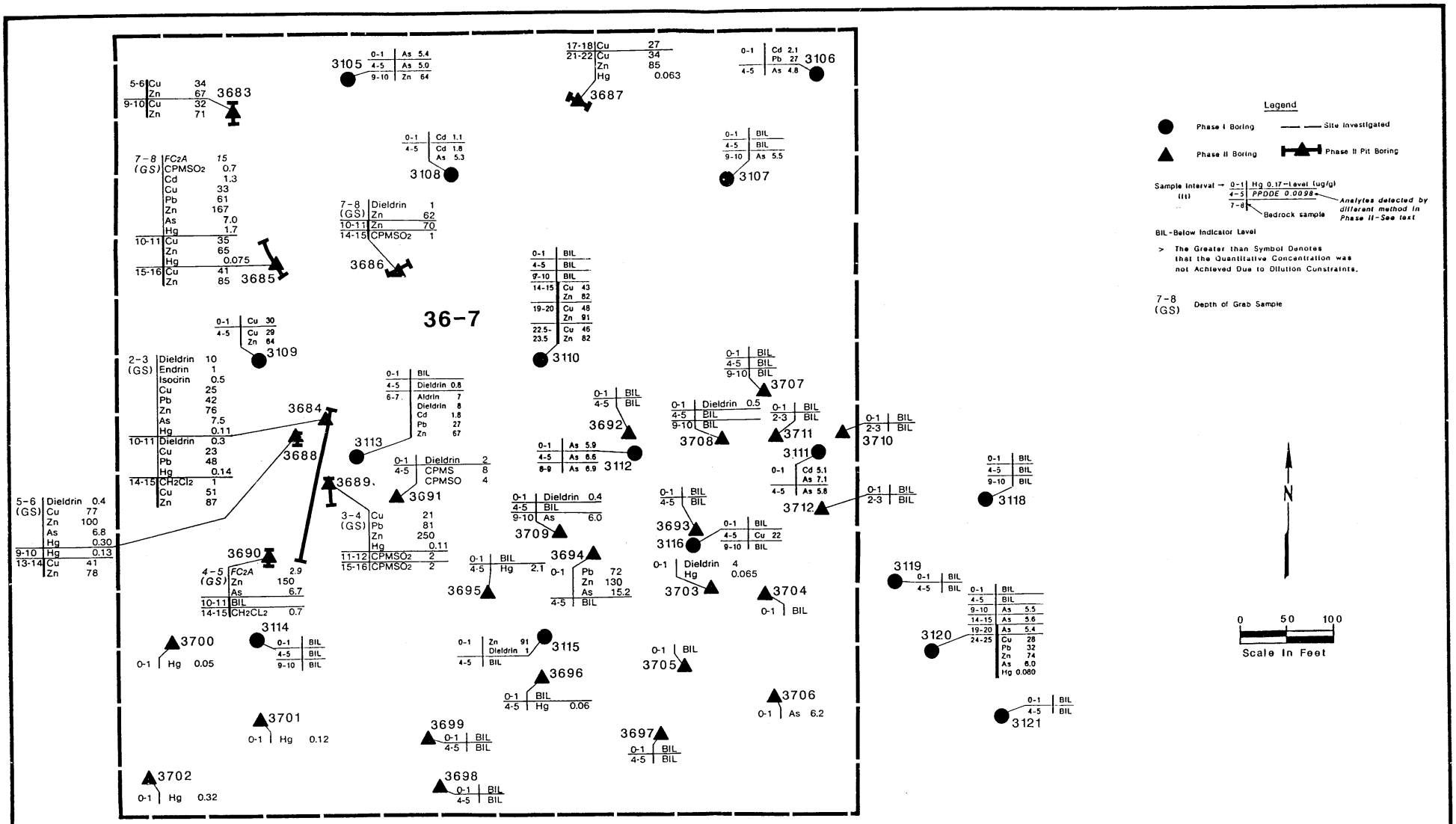


Figure 36-7-II-2

SITE 36-7, PHASE I AND PHASE II INVESTIGATIONS
CHEMICAL ANALYSIS RESULTS

SOURCE: HLA, 1988

Prepared for:
U.S. Army Program Manager's Office
For Rocky Mountain Arsenal

Aberdeen Proving Ground, Maryland

Dieldrin was detected in the 0- to 1-ft sample interval from Boring 3703 at a concentration of 4 ppm. This boring was located adjacent to the northwest corner of the Shell incinerator. Dieldrin was also detected in Boring 3708 (0.5 ppm, 0- to 1-ft interval) and Boring 3709 (0.4 ppm, 0- to 1-ft interval). Boring 3708 is located at the intersection of the entrance to the incinerator, and Boring 3709 is in the fill area within Anomaly G.

The organosulphur compounds CPMS and CPMSO were detected at concentrations of 8 ppm and 4 ppm, respectively, in the 4- to 5-ft sample interval of Boring 3691. Four samples obtained at depths greater than 7 ft contained $CPMSO_2$ at concentrations ranging from 0.7 to 2 ppm. Three of these samples were collected below the base of suspected disposal trenches.

Of the ICP metals, cadmium, copper, lead, and zinc were detected in 1, 14, 5, and 15 samples, respectively, at concentrations within or above their indicator ranges. High concentrations of zinc, copper, and lead were detected in samples obtained from four pit borings (3684, 3688, 3689, and 3690) located in Anomaly B. With the exception of one lead and one zinc value, all of the ICP metal values within or above indicator ranges are associated with the grab samples of trench material or bedrock samples from the pit borings.

Seven of the 56 samples analyzed for arsenic were within or above the indicator range. Arsenic concentrations were in the middle of the indicator range, except Boring 3694 (0 to 1 ft) which contained arsenic at 15 ppm. Mercury was detected in 14 of 56 samples at concentrations ranging from 0.054 to 2.1 ppm. Higher mercury values are generally associated with samples obtained from the pit borings.

Grab samples from various depth intervals (see Table 36-7-II-3) were analyzed for TDGCL, IMPA, FC2A, and MPA. FC2A was detected in grab samples obtained from Pit Boring 3685 (7- to 8-ft interval) and Pit Boring 3690 (4- to 5-ft interval) at concentrations of 15.2 and 2.9 ppm, respectively.

The data reporting procedures as described in the Laboratory Quality Assurance Plan for RMA (ESE, 1985, Appendix B, RIC#85127R07) require that all analyses on a sample be completed within their respective holding time and that analytical results be corrected for percent recovery and moisture content. During routine sample analysis, analytical results must either fall within or be diluted within the Certified Range provided that holding times have not expired.

During laboratory certification, an analytical method is tested over a certain concentration range to determine the Certified Range. A typical tested concentration range would be 0, 0.5X, 1.0X, 2.0X, 5.0X, and 10.0X, where X is the Target Reporting Limit (TRL). The Certified Reporting Limit (CRL) is determined by comparing the target and actual concentrations of the tested range. The upper Certified Range is the higher target concentration achieved.

If a sample analysis indicates that the sample was not diluted adequately to be within the Certified Range, the result is reported as greater than (>) the upper Certified Range times any dilution factors. If a sample has exceeded its holding time and the result is greater than the Certified Range, the result is reported as greater than the upper Certified Range. If holding times are exceeded in attempting to dilute the sample until all results are within the Certified Range, results that are not identified above the Certified Range but that may be present at concentrations above the certified detection limit are reported as the detection limit times the dilution factor.

Several compounds detected by GC/MS were not included in the target compound list and were not conclusively identified. These compounds are included in the data presented in Appendix 36-7-II-B. Table 36-7-II-4 summarizes nontarget compounds detected at Site 36-7. It should be noted that an individual compound may have more than one retention time and that a particular retention time may be assigned to more than one compound. Table 36-7-II-4, therefore, provides only a general indication of additional compounds that may be present.

Table 36-7-II-4. Tentative Identification of Nontarget Compounds in Site 36-7 Phase II Soil Samples (page 1 of 4)

Boring Number	Interval Depth (ft)	Unknown Number	Concentration Above Background (ppm)*	Sample Number	Lot	Best Fit	Comments
3683	5-6	526	10	36-7-47X2	KXI	Methylhexanone	
		614	3			Dibutyl nonanedioate	d
	9-10	526	10	36-7-47X3	KXI	Methylhexanone	
		614	4			Dibutyl nonanedioate	d
3684	2-3	587	0.9	36-7-47X5	KES	Unknown	a,f
		588	1			Unknown	a
		593	2			Acetylhydroxydimethylpyridinone	
	10-11	524	8	36-7-47X6	KXG	Methylhexanone	
		549	0.9			Pentanoic acid	d,f
		554	1			Unknown	a
		558	1			Unknown	a
		566	1			Nonanoic acid	d
		586	2			Dodecanoic acid	d
		598	2			Tetradecanoic acid	d
	14-15	160	9	36-7-47X7	KZQ	Hexamethylcyclotrisiloxane	i
		524	8		KXG	Methylhexanone	
		558	0.9			Unknown	a,f
		566	1			Nonanoic acid	d
		586	1			Dodecanoic acid	d
		598	1			Tetradecanoic acid	d
		615	2			Dibutyl nonanedioate	d
3685	7-8	543	1	36-7-47X9	KES	Unknown	a
		559	2			Octanoic acid	d
		566	0.9			Nonanoic acid	d,f
		573	0.9			Decanoic acid	d,f
		576	10			Trichlorobenzenamine	
		578	10			Unknown	a
		588	2			Unknown	a
		598	3			2-Chlorophosphate ethanol (3:1)	
		615	2			Unknown alkane	a
		622	1			Unknown	a
		624	0.9			Unknown alkane	a,f
		626	1			Unknown	a
		628	4			Dioctyl hexanedioate	d
		631	4			Unknown	a
		632	2			Unknown alkane	a
		634	1			Bis (2-ethylhexyl) phthalate	c
		636	1			Unknown alkane	a
		640	1			Unknown alkane	a
		644	1			Unknown alkane	a
		649	2			Unknown alkane	a
		655	2			Unknown alkane	a

Table 36-7-II-4. Tentative Identification of Nontarget Compounds in Site 36-7 Phase II Soil Samples (page 2 of 4)

Boring Number	Interval Depth (ft)	Unknown Number	Concentration Above Background (ppm)*	Sample Number	Lot	Best Fit	Comments
	10-11	526	10	36-7-47X10	KXI	Methylhexanone	
		614	2			Dibutyl nonanedioate	d
	15-16	526	10	36-7-47X12	KXI	Methylhexanone	
		614	3			Dibutyl nonanedioate	d
3686	6-7	576	2	36-7-47X13	KET	Trichlorobenzenamine	
		577	2			Junipene	
		594	0.8			Unknown alkane	a,f
		602	1			Phenylethylphenol	
		629	0.9			Unknown	a,f
		630	7			Methylsulfonyldinitrodipropylbenzenamine	
		636	1			Unknown	a
		637	3			Unknown	a
	10-11	526	10	36-7-47X14	KXI	Methylhexanone	
		614	3			Dibutyl nonanedioate	d
	14-15	160	3	36-7-47X15	KZO	Hexamethylcyclotrisiloxane	i
		526	10		KXI	Methylhexanone	
3687	17-18	526	20	36-7-47X18	KXI	Methylhexanone	
	21-22	160	0.8	36-7-47X19	KZO	Hexamethylcyclotrisiloxane	i,f
		526	10		KXI	Methylhexanone	
		555	2			Unknown	a
		556	1			Unknown	a
3688	5-6	600	0.8	36-7-47X17	KET	Unknown alkane	a,f
		630	0.8			2-Ethylhexyldiphenyl phosphonate	f
	9-10	550	8	36-7-47X103	KXG	Unknown	a
		554	6			Unknown	a
		572	40			Trichlorophenol	
		626	10			Ethylhexyl trichlorophenoxyethanoate	
		627	400			Ethylhexyl trichlorophenoxyethanoate	
		628	40			Ethylhexyl trichlorophenoxyethanoate	
		629	20			Ethylhexyl trichlorophenoxyethanoate	
		630	200			Ethylhexyl trichlorophenoxyethanoate	
		631	600			Ethylhexyl trichlorophenoxyethanoate	
		632	100			Ethylhexyl trichlorophenoxyethanoate	
		633	20			Ethylhexyl trichlorophenoxyethanoate	
		634	20			Ethylhexyl trichlorophenoxyethanoate	
	13-14	160	6	36-7-47X104	KZQ	Hexamethylcyclotrisiloxane	i
		524	6		KXG	Methylhexanone	
		566	1			Nonanoic acid	d
		586	1			Dodecanoic acid	d
		614	0.9			Dibutyl nonanedioate	d,f

Table 36-7-II-4. Tentative Identification of Nontarget Compounds in Site 36-7 Phase II Soil Samples (page 3 of 4)

Boring Number	Interval Depth (ft)	Unknown Number	Concentration Above Background (ppm)*	Sample Number	Lot	Best Fit	Comments
3689	3-4	582	4	36-7-47X21	KET	Dihydroacenaphthylene	
		598	1			Unknown	a
		617	2			Pyrene	
		618	0.8			Phenylnaphthalene	f
		624	1			Unknown alkane	a
		628	1			Unknown alkane	a
		632	1			Unknown alkane	a
		633	0.9			Triphenylene	f
		636	1			Unknown alkane	a
		640	1			Unknown alkane	a
	11-12	524	10	36-7-47X22	KXG	Methyl hexanone	
		632	2			Methylsulfonyldinitrodipropylbenzenamine	
	15-16	161	3	36-7-47X23	KZQ	Hexamethyltrisiloxane	i
		524	9		KXG	Methylhexanone	
		541	1			Unknown	a
		614	0.9			Dibutyl nonanedioate	d,f
		636	3			Bis (2-ethylhexyl) phthalate	c
3690	4-5	551	0.9	36-7-47X25	KET	Heptanoic acid	d,f
		576	7			Trichlorobenzenenmamine	
		578	6			Unknown	a
		591	0.9			Unknown alkane	a,f
		594	2			Unknown alkane	a
		595	3			Unknown alkane	a
		600	2			Unknown alkane	a
	10-11	524	9	36-7-47X26	KXG	Methylhexanone	
	14-15	524	9	36-7-47X27	KXG	Methylhexanone	
		615	0.9			Dibutylnonanedioate	d
3691	0-1	526	9	36-7-47X29	KXI	Methylhexanone	
3694	0-1	518	1	36-7-47X41	KXG	Tetrachloroethene	
		524	8			Methylhexanone	
		566	0.8			Nonanoic acid	d,f
		586	0.7			Dodecanoic acid	d,f
		661	0.8			Unknown alkane	a,f
	4-5	524	8	36-7-47X42	KXG	Methylhexanone	
3696	0-1	526	10	36-7-47X49	KXI	Methylhexanone	
		619	0.9			Unknown	a,f
		628	0.9			Butyl octadecanoate	d,f
		633	0.8			Unknown	a,f
	4-5	526	10	36-7-47X50	KXI	Methylhexanone	
3697	0-1	635	0.8	36-7-47X53	KXI	Bis (2-ethylhexyl) phthalate	c,f
	4-5	635	2	36-7-47X54	KXI	Bis (2-ethylhexyl) phthalate	c

Table 36-7-II-4. Tentative Identification of Nontarget Compounds in Site 36-7 Phase II Soil Samples (page 4 of 4)

Boring Number	Interval Depth (ft)	Unknown Number	Concentration Above Background (ppm)*	Sample Number	Lot	Best Fit	Comments
3698	0-1	524	9	36-7-47X57	KXG	Methylhexanone	
		558	0.8			Unknown	a,f
		566	1			Nonanoic acid	d
		586	0.9			Dodecanoic acid	d,f
	4-5	524	8	36-7-47X58	KXG	Methylhexanone	
3700	0-1	636	1	36-7-47X65	KXC	Bis (2-ethylhexyl) phthalate	c
3703	0-1	523	0.9	36-7-47X71	KXC	Oxabicycloheptane	f
		616	3			Unknown	a
		632	2			Methylsulfonyldinitrodipropylbenzenamine	
		673	5			Unknown	a
3704	0-1	636	2	36-7-47X73	KXC	Bis (2-ethylhexyl) phthalate	c
3705	0-1	637	0.9	36-7-47X75	KXC	Bis (2-ethylhexyl) phthalate	c,f
3706	0-1	615	1	36-7-47X77	KXC	Dibutyl nonanedioate	d
		636	1			Bis (2-ethylhexyl) phthalate	c
3707	0-1	524	1	36-7-47X79	KXC	Methylhexanone	
	9-10	160	1	36-7-47X81	KZN	Hexamethyltrisiloxane	i
		625	0.9		KXF	Unknown phthalate	a,c,f
		628	0.8			Unknown phthalate	a,c,f
		636	4			Bis (2-ethylhexyl) phthalate	c
3708	4-5	636	0.8	36-7-47X85	KXF	Bis (2-ethylhexyl) phthalate	c,f
	9-10	160	6	36-7-47X86	KZN	Hexamethyltrisiloxane	i
		636	1		KXF	Bis (2-ethylhexyl) phthalate	c
3709	0-1	630	3	36-7-47X89	KXE	Bis (2-ethylhexyl) phthalate	c
		636	8			Dioctyl hexanedioate	d
	9-10	615	0.9	36-7-47X91	KXE	Dibutyl nonanedioate	d,f

* Values reported are method blank corrected.

\+ a. No positive identification.
b. Surfactant.
c. Plasticizer (Note: All phthalates and adipates will have this comment).
d. Derived from natural products.
e. Suspected laboratory contaminant.
f. Low concentration.
g. Low frequency of occurrence.
h. Ubiquitous.
i. Possible column bleed.
j. None detected.

Source: ESE, 1988.

Nontarget compounds were detected in 42 of the 56 samples analyzed by GC/MS. Methyl hexanone was tentatively identified in 23 samples from Lots KXI, KXG, and KXC at concentrations ranging from 1 to 20 ppm. Hexamethyltrisiloxane, which was identified in 7 of 11 samples analyzed by GC/MS for VO compounds, is associated with column bleed during laboratory procedures.

Trichlorobenzenamine and methylsulfonyldinitrodipropylbenzenamine were each tentatively identified in three samples. The grab sample (6 to 7 ft) from Boring 3686 contained phenyethylphenol (2 ppm) and junipene (1 ppm) in addition to trichlorobenzene and methylsulfonyldinitrodipropylbenzenamine. Boring 3688 (9 to 10 ft) contained trichlorophenol and ethylhexyl trichlorophenoxy-ethanoate at elevated concentrations ranging from 10 to 600 ppm. Pyrene, triphenylene, and two naphthalenes were discovered in Boring 3689 (3 to 4 ft). The remaining comounds were predominantly naturally occurring compounds, phthalates, or could not be conclusively identified.

Results of the Phase II sampling program at Site 36-7 will be included as part of the overall analysis of the Central Study Area Report.

5.0 REFERENCES

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APPENDIX 36-7-II-A
CHEMICAL NAMES, METHODS, AND ABBREVIATIONS

APPENDIX 36-7-II-A
CHEMICAL NAMES, METHODS, AND ABBREVIATIONS

PHASE I ANALYTES AND CERTIFIED METHODS

Analytes/Methods	Synonymous Names and Abbreviations	Standard Abbreviations
VOLATILE ORGANIC COMPOUNDS/GCMS	VOL	VO
1,1-Dichloroethane	1,1-Dichloroethane	11DCLE
1,2-Dichloroethane	1,2-Dichloroethane	12DCLE
1,1,1-Trichloroethane (TCA)	1,1,1-Trichloroethane	111TCE
1,1,2-Trichloroethane	1,1,2-Trichloroethane	112TCE
Benzene	Benzene	C_6H_6
Bicycloheptadiene	Bicycloheptadiene (BCHD)	BCHPD
Carbon tetrachloride	Carbon tetrachloride	CCL_4
Chlorobenzene	Chlorobenzene	CLC_6H_5
Chloroform	Chloroform	$CHCL_3$
Dibromochloropropane	Dibromochloropropane	DBCP
Dicyclopentadiene	Dicyclopentadiene	DCPD
Dimethyldisulfide	Dimethyldisulfide	DMDS
Ethylbenzene	Ethylbenzene	ETC_6H_5
m-Xylene	meta-Xylene	13DMB
Methylene chloride	Methylene chloride	CH_2CL_2
Methylisobutyl ketone	Methylisobutyl ketone	MIBK
o,p-Xylene	ortho- and/or para-Xylene	XYLEN
Tetrachloroethene (PCE)	Tetrachloroethylene	TCLEE
Toluene	Toluene	MEC_6H_5
Trans 1,2-dichloroethene	Trans 1,2-dichloroethylene	12DCE
Trichloroethene (TCE)	Trichloroethylene	TRCLE
SEMIVOLATILE ORGANIC COMPOUNDS/GCMS	EXTRACTABLE ORGANIC COMPOUNDS (EX)	SVO
1,4-Oxathiane	1,4-Oxathiane	OXAT
2,2-Bis (para-chlorophenyl)-1,1-dichloroethane	Dichlorodiphenylethane	PPDDE
2,2-Bis (para-chlorophenyl) 1,1,1-trichloroethane	Dichlorodiphenyltrichloroethane	PPDDT
Aldrin	Aldrin	ALDRN
Atrazine	Atrazine	ATZ
Chlordane	Chlordane	CLDAN
Chlorophenylmethyl sulfide	p-Chlorophenylmethyl sulfide	CPMS
Chlorophenylmethyl sulfoxide	p-Chlorophenylmethyl sulfoxide	CPMSO
Chlorophenylmethyl sulfone	p-Chlorophenylmethyl sulfone	$CPMSO_2$
Dibromochloropropane	Dibromochloropropane	DBCP
Dicyclopentadiene	Dicyclopentadiene	DCPD
Dieldrin	Dieldrin	DLDRN
Diisopropylmethyl phosphonate	Diisopropylmethyl phosphonate	DIMP

APPENDIX 36-7-II-A
CHEMICAL NAMES, METHODS, AND ABBREVIATIONS

Analytes/Methods	Synonymous Names and Abbreviations	Standard Abbreviations
SEMIVOLATILE ORGANIC COMPOUNDS (CONT)		
Dimethylmethyl phosphonate	Dimethylmethyl phosphonate	DMMP
Dithiane	Dithiane	DITH
Endrin	Endrin	ENDRN
Hexachlorocyclopentadiene	Hexachlorocyclopentadiene (HCPD)	CL_6CP
Isodrin	Isodrin	ISODR
Malathion	Malathion	MLTHN
Parathion	Parathion	PRTHN
Supona	2-Chloro-1(2,4-dichlorophenyl) vinyldiethyl phosphate	SUPONA
Vapona	Vapona	DDVP
METALS/ICP	ICAP	ICP
Cadmium	Cadmium	CD
Chromium	Chromium	CR
Copper	Copper	CU
Lead	Lead	PB
Zinc	Zinc	ZN
SEPARATE ANALYSES		
Arsenic/AA	Arsenic	AS
Mercury/AA	Mercury	HG
Dibromochloropropane/GC	Dibromochloropropane	DBCP

APPENDIX 36-7-II-A
CHEMICAL NAMES, METHODS, AND ABBREVIATIONS

PHASE II ANALYTES AND CERTIFIED METHODS

Analytes/Methods	Synonymous Names and Abbreviations	Standard Abbreviations
VOLATILE ORGANIC COMPOUNDS/GCMS (Same as Phase I)	VOL	VO
SEMIVOLATILE ORGANIC COMPOUNDS/GCMS (Same as Phase I)	EXTRACTABLE ORGANIC COMPOUNDS (EX)	SVO
VOLATILE HALOCARBON COMPOUNDS/GCCON	PURGEABLE HALOCARBONS (PHC)	VHO
1,1-Dichloroethane	1,1-Dichloroethane	11DCLE
1,2-Dichloroethane	1,2-Dichloroethane	12DCLE
1,1-Dichloroethene	1,1-Dichloroethene	11DCE
1,1,1-Trichloroethane (TCA)	1,1,1-Trichloroethane	111TCE
1,1,2-Trichloroethane	1,1,2-Trichloroethane	112TCE
Carbon tetrachloride	Carbon tetrachloride	CCL_4
Chlorobenzene	Chlorobenzene	CLC_6H_5
Chloroform	Chloroform	$CHCL_3$
Methylene chloride	Methylene chloride	CH_2CL_2
Trans 1,2-dichloroethylene	Trans 1,2-dichloroethene	12DCE
Tetrachloroethene (PCE)	Tetrachloroethylene	TCLEE
Trichloroethene (TCE)	Trichloroethylene	TRCLE
VOLATILE HYDROCARBON COMPOUNDS/GCFID	DCPD	HYDCBN
Bicycloheptadiene	Bicycloheptadiene (BCHD)	BCHPD
Dicyclopentadiene	Dicyclopentadiene	DCPD
Methylisobutyl ketone	Methylisobutyl ketone	MIBK
VOLATILE AROMATIC COMPOUNDS/GCPID	PURGEABLE AROMATICS (PAM)	VAO
Benzene	Benzene	C_6H_6
Ethylbenzene	Ethylbenzene	ETC_6H_5
m-Xylene	meta-Xylene	13DMB
o,p-Xylene	ortho- and/or para-Xylene	XYLEN
Toluene	Toluene	MEC_6H_5
ORGANOCHLORINE PESTICIDES/GCEC		OCP
2,2-Bis (para-chlorophenyl)-1,1-dichloroethane	Dichlorodiphenylethane	PPDDE
2,2-Bis (para-chlorophenyl)-1,1,1-trichloreoethane	Dichlorodiphenyltrichloroethane	PPDDT
Aldrin	Aldrin	ALDRN
Chlordane	Chlordane	CLDAN
Dieldrin	Dieldrin	DLDRN
Endrin	Endrin	ENDRN
Hexachlorocyclopentadiene	Hexachlorocyclopentadiene	CL_6CP
Isodrin	Isodrin	ISODR

APPENDIX 36-7-II-A
CHEMICAL NAMES, METHODS, AND ABBREVIATIONS

Analytes/Methods	Synonymous Names and Abbreviations	Standard Abbreviations
ORGANOPHOSPHOROUS PESTICIDES/GCNPD	ORGANOPHOSPHOROUS COMPOUNDS (OPC)	OPP
Atrazine	Atrazine	ATZ
Malathion	Malathion	MLTHN
Parathion	Parathion	PRTHN
Supona	2-Chloro-1(2,4-dichlorophenyl) vinyldiethyl phosphate	SUPONA
Vapona	Vapona	DDVP
ORGANOPHOSPHOROUS COMPOUNDS/GCFPD	DIMP	OPC
Diisopropylmethyl phosphonate	Diisopropylmethyl phosphonate	DIMP
Dimethylmethyl phosphonate	Dimethylmethyl phosphonate	DMMP
ORGANOSULPHUR COMPOUNDS/GCFPD		OSC
1,4-Oxathiane	1,4-Oxathiane	OXAT
Benzothiazole	Benzothiazole	BTZ
Chlorophenylmethyl sulfide	p-Chlorophenylmethyl sulfide	CPMS
Chlorophenylmethyl sulfone	p-Chlorophenylmethyl sulfone	$CPMSO_2$
Chlorophenylmethyl sulfoxide	p-Chlorophenylmethyl sulfoxide	CPMSO
Dimethyldisulfide	Dimethyldisulfide	DMDS
Dithiane	Dithiane	DITH
METALS/ICP	ICAP	ICP
Cadmium	Cadmium	CD
Chromium	Chromium	CR
Copper	Copper	CU
Lead	Lead	PB
Zinc	Zinc	ZN
SEPARATE ANALYSES		
Arsenic/AA	Arsenic	AS
Mercury/AA	Mercury	HG
Dibromochloropropane/GC	Dibromochloropropane	DBCP

APPENDIX 36-7-II-A
CHEMICAL NAMES, METHODS, AND ABBREVIATIONS

Analytes/Methods	Synonymous Names and Abbreviations	Standard Abbreviations
ARMY AGENT DEGRADATION PRODUCTS:		ADP
AGENT PRODUCTS/HPLC	TDGCL	
Chloroacetic Acid	Chloroacetic acid	CLC2A
Thiodiglycol	Thiodiglycol (TDG)	TDGCL
AGENT PRODUCTS/IONCHROM	IMPA	GBDP
Fluoroacetic acid	Fluoroacetic acid	FC2A
Isopropylmethylphosphonic acid	Isopropylmethylphosphonate	IMPA
Methylphosphonic acid	Methylphosphonate	MPA

Methods	Abbreviations
Atomic Absorption Spectroscopy	AA
Gas Chromatography/Conductivity Detector	GCCON
Gas Chromatography/Electron Capture	GCEC
Gas Chromatography/Flame Ionization Detector	GCFID
Gas Chromatography/Flame Photometric	GCFPD
Gas Chromatography/Mass Spectrometry	GCMS
Gas Chromatography/Nitrogen Phosphorous Detector	GCNPD
Gas Chromatography/Photoionizaton Detector	GCPID
High Performance Liquid Chromatography	HPLC
Inductively Coupled Argon Plasma	ICP, ICAP
Ion Chromatography	IONCHROM

APPENDIX 36-7-II-B
PHASE II CHEMICAL DATA

ENVIRONMENTAL SCIENCE & ENGINEERING 06/24/88 PAGE#1
PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	STORET #	SAMPLE ID/# 3683A 36-7-47 2	3683B 36-7-47 3	3684GRAB 36-7-47 5	3684A 36-7-47 6	3684B 36-7-47 7	3685GRAB 36-7-47 9	3685A 36-7-47 10	3685X 36-7-47 12	3686GRAB 36-7-47 13	3686A 36-7-47 14	3686B 36-7-47 15	3688GRAB 36-7-47 17	3687A 36-7-47 18	3687B 36-7-47 19	3689GRAB 36-7-47 21
UNITS	METHOD															
DATE		03/15/88	03/15/88	02/03/88	03/16/88	03/16/88	02/02/88	03/15/88	03/15/88	02/02/88	03/14/88	03/14/88	02/03/88	03/14/88	03/14/88	02/03/88
TIME		08:19	08:33	12:10	08:43	08:54	12:22	09:20	10:12	10:57	09:54	10:22	12:48	09:40	12:03	10:16
SAMPLE TYPE	71999 0	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO
SITE TYPE 1	99759 0	BORE	BORE	CSDT	BORE	BORE	CSDT	BORE	BORE	CSDT	BORE	BORE	CSDT	BORE	BORE	CSDT
SAMPLE DEPTH FT	99758 0	5	9	0	10	14	0	10	15	0	10	14	0	17	21	0
SAMPLING TECHNIQUE	72005 0	S	S	G	S	S	S	S	S	G	S	S	G	S	S	G
INSTALLATION CODE SAMPLE	99720 0	RK	RK	RK	RK	RK	RKO	RK	RK	RK	RK	RK	RK	RK	RK	RK
MOISTURE %WET WT	70320 I	24.1	23.8	12.1	13.5	21.7	19.6	25.4	22.2	13.3	15.7	20.7	13.0	24.2	21.9	10.8
CADMIUM UG/G- DRY	1028 R9	<0.921	<0.921	<0.921	<0.921	<0.921	1.28	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921
CHROMIUM UG/G-DRY	99584 R9	<7.16	<7.16	16.9	9.48	<7.16	11.8	<7.16	<7.16	11.5	<7.16	<7.16	15.0	<7.16	<7.16	<7.16
COPPER UG/G- DRY	1043 R9	34.3	31.6	24.7	22.9	51.3	32.9	35.0	41.0	17.7	12.3	19.6	77.1	27.0	34.4	21.4
LEAD UG/G-DRY	1052 R9	<16.8	<16.8	41.8	48.0	<16.8	61.4	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	80.6
ZINC UG/G-DRY	1093 R9	66.6	70.5	75.6	48.8	87.3	167	64.9	84.7	61.8	70.1	54.7	100	58.2	84.6	252
ARSENIC UG/G- DRY	1003 T9	<4.70	<4.70	7.45	<4.70	<4.70	6.97	<4.70	<4.70	<4.70	<4.70	<4.70	6.83	<4.70	<4.70	<4.70
MERCURY UG/G-DRY	71921 V9	<0.050	<0.050	0.110	0.138	<0.050	1.67	0.075	<0.050	<0.050	<0.050	<0.050	0.303	<0.050	0.063	0.105
ALDRIN UG/G-DRY	98356 Q9	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94
ATRAZINE UG/G-DRY	98655 Q9	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73
CHLORDANE UG/G-DRY	98361 Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
P-CLPHENYLMETHY-SULFIDE UG/G-DRY	98653 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
P-CLPHENYLMETHY-SULFOXIDE UG/G-DRY	98654 Q9	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35
P-CLPHENYLMETHY-SULFONE UG/G-DRY	98703 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	0.71	<0.29	<0.29	<0.29	<0.29	1.3	<0.29	<0.29	<0.29	<0.29
DBCP(NEMAGON) UG/G-DRY	98652 Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
DICYCLOPENTADIENE UG/G-DRY	98651 Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
DDE,PP' UG/G-DRY	98363 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
DDT,PP' UG/G-DRY	98364 Q9	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37
DIELDRIN UG/G-DRY	98365 Q9	<0.25	<0.25	9.8	0.33	<0.25	<0.25	<0.25	<0.25	1.2	<0.25	<0.25	0.35	<0.25	<0.25	<0.25

PARAMETERS	UNITS	STORET #	METHOD	SAMPLE ID/# 3683A 36-7-47 2	3683B 36-7-47 3	3684GRAB 36-7-47 5	3684A 36-7-47 6	3684B 36-7-47 7	3685GRAB 36-7-47 9	3685A 36-7-47 10	3685X 36-7-47 12	3686GRAB 36-7-47 13	3686A 36-7-47 14	3686B 36-7-47 15	3688GRAB 36-7-47 17	3687A 36-7-47 18	3687B 36-7-47 19	3689GRAB 36-7-47 21
DATE				03/15/88	03/15/88	02/03/88	03/16/88	03/16/88	02/02/88	03/15/88	03/15/88	02/02/88	03/14/88	03/14/88	02/03/88	03/14/88	03/14/88	02/03/88
TIME				08:19	08:33	12:10	08:43	08:54	12:22	09:20	10:12	10:57	09:54	10:22	12:48	09:40	12:03	10:16
DIMP	UG/G-DRY	98645	Q9	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,4 DITHIANE	UG/G-DRY	98650	Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
DMMP	UG/G-DRY	98657	Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
ENDRIN	UG/G-DRY.	98369	Q9	<0.70	<0.70	1.4	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70
HEXACHLOROCYCLOPENTADIENE	UG/G-DRY	98647	Q9	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1
ISODRIN	UG/G-DRY	98649	Q9	<0.33	<0.33	0.53	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
MALATHION	UG/G-DRY	98648	Q9	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59
1,4 OXATHIANE	UG/G-DRY	98644	Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
ETY'PARATHION	UG/G-DRY	98658	Q9	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63
SUPONA	UG/G-DRY	98656	Q9	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49
VAPONA	UG/G-DRY	98646	Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
DICYCLOPENTADIENE	UG/G-DRY	98651	W9		<0.27			<0.27			<0.27			<0.27			<0.27	
TRANS-1,2-DICHLOROETHENE	UG/G-DRY	98687	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
ETHYLBENZENE	UG/G-DRY	98688	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
METHYLENE CHLORIDE	UG/G-DRY	98689	W9		<0.25			0.97			<0.25			<0.25			<0.25	
TETRACHLOROETHENE	UG/G-DRY	98690	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
TOLUENE	UG/G-DRY	98691	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
1,1,1-TRICHLOROETHANE	UG/G-DRY	98692	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
1,1,2-TRICHLOROETHANE	UG/G-DRY	98693	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
TRICHLOROETHENE	UG/G-DRY	98694	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
M-XYLENE	UG/G-DRY	98695	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
MIBK	UG/G-DRY	98696	W9		<0.50			<0.50			<0.50			<0.50			<0.50	
DMDS	UG/G-DRY	98697	W9		<0.25			<0.25			<0.25			<0.25			<0.25	
BENZENE	UG/G-DRY	98699	W9		<0.25			<0.25			<0.25			<0.25			<0.25	

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PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3683A 36-7-47 2	3683B 36-7-47 3	3684GRAB 36-7-47 5	3684A 36-7-47 6	3684B 36-7-47 7	3685GRAB 36-7-47 9	3685A 36-7-47 10	3685X 36-7-47 12	3686GRAB 36-7-47 13	3686A 36-7-47 14	3686B 36-7-47 15	3688GRAB 36-7-47 17	3687A 36-7-47 18	3687B 36-7-47 19	3689GRAB 36-7-47 21
DATE			03/15/88	03/15/88	02/03/88	03/16/88	03/16/88	02/02/88	03/15/88	03/15/88	02/02/88	03/14/88	03/14/88	02/03/88	03/14 88	03/14/88	02/03/88
TIME			08:19	08:33	12:10	08:43	08:54	12:22	09:20	10:12	10:57	09:54	10:22	12:48	09:40	12:03	10:16
UNK600		90600				0.9								0.8			
	UG/G	Q9															
UNK609		90609				5											
	UG/G	Q9															
UNK618		90618				2											0.8
	UG/G	Q9															
UNK619		90619				1											
	UG/G	Q9															
UNK626		90626				0.8		1									
	UG/G	Q9															
UNK631		90631				0.9		4									
	UG/G	Q9															
UNK633		90633				2											0.9
	UG/G	Q9															
UNK160		90160					9						3			0.8	
	UG/G	W9															
UNK615		90615					2	2									
	UG/G	Q9															
UNK543		90543						1									
	UG/G	Q9															
UNK559		90559						2									
	UG/G	Q9															
UNK573		90573						0.9									
	UG/G	Q9															
UNK576		90576						10			2						
	UG/G	Q9															
UNK578		90578						10									
	UG/G	Q9															
UNK622		90622						1									
	UG/G	Q9															
UNK624		90624						0.9									1.0
	UG/G	Q9															
UNK628		90628						4									1
	UG/G	Q9															
UNK632		90632						2									1
	UG/G	Q9															
UNK634		90634						1									
	UG/G	Q9															
UNK636		90636						1			1						1
	UG/G	Q9															
UNK640		90640						1.0									1
	UG/G	Q9															
UNK644		90644						1.0									
	UG/G	Q9															
UNK649		90649						2									
	UG/G	Q9															
UNK655		90655						2									
	UG/G	Q9															

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3683A 36-7-47 2	3683B 36-7-47 3	3684GRAB 36-7-47 5	3684A 36-7-47 6	3684B 36-7-47 7	3685GRAB 36-7-47 9	3685A 36-7-47 10	3685X 36-7-47 12	3686GRAB 36-7-47 13	3686A 36-7-47 14	3686B 36-7-47 15	3688GRAB 36-7-47 17	3687A 36-7-47 18	3687B 36-7-47 19	3689GRAB 36-7-47 21
DATE			03/15/88	03/15/88	02/03/88	03/16/88	03/16/88	02/02/88	03/15/88	03/15/88	02/02/88	03/14/88	03/14/88	02/03/88	03/14/88	03/14/88	02/03/88
TIME			08:19	08:33	12:10	08:43	08:54	12:22	09:20	10:12	10:57	09:54	10:22	12:48	09:40	12:03	10:16
UNK577		90577									2						
	UG/G	Q9															
UNK594		90594									0.8						
	UG/G	Q9															
UNK629		90629									0.9						
	UG/G	Q9															
UNK630		90630									7			0.8			
	UG/G	Q9															
UNK637		90637									3						
	UG/G	Q9															
UNK555		90555														2	
	UG/G	Q9															
UNK556		90556														1	
	UG/G	Q9															
UNK582		90582															4
	UG/G	Q9															
UNK617		90617															2
	UG/G	Q9															
UNK161		90161															
	UG/G	W9															
UNK541		90541															
	UG/G	Q9															
UNK551		90551															
	UG/G	Q9															
UNK591		90591															
	UG/G	Q9															
UNK595		90595															
	UG/G	Q9															
UNK518		90518															
	UG/G	Q9															
UNK661		90661															
	UG/G	Q9															
UNK523		90523															
	UG/G	Q9															
UNK616		90616															
	UG/G	Q9															
UNK673		90673															
	UG/G	Q9															
UNK550		90550															
	UG/G	Q9															
UNK572		90572															
	UG/G	Q9															
UNK627		90627															
	UG/G	Q9															
UNK526		90526	10	10					10	10		10	10		20	10	
	UG/G	Q9															
UNK614		90614	3	4	1				2	3		3					1
	UG/G	Q9															

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PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1

FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER

ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3683A 36-7-47 2	3683B 36-7-47 3	3684GRAB 36-7-47 5	3684A 36-7-47 6	3684B 36-7-47 7	3685GRAB 36-7-47 9	3685A 36-7-47 10	3685X 36-7-47 12	3686GRAB 36-7-47 13	3686A 36-7-47 14	3686B 36-7-47 15	3688GRAB 36-7-47 17	3687A 36-7-47 18	3687B 36-7-47 19	3689GRAB 36-7-47 21
DATE			03/15/88	03/15/88	02/03/88	03/16/88	03/16/88	02/02/88	03/15/88	03/15/88	02/02/88	03/14/88	03/14/88	02/03/88	03/14/88	03/14/88	02/03/88
TIME			08:19	08:33	12:10	08:43	08:54	12:22	09:20	10:12	10:57	09:54	10:22	12:48	09:40	12:03	10:16
UNK635	UG/G	90635 Q9		0.2													

PARAMETERS	UNITS	STORET #	METHOD	3689GRAB 36-7-47 21	3689A 36-7-47 22	3689B 36-7-47 23	3690GRAB 36-7-47 25	3690A 36-7-47 26	3690B 36-7-47 27	3691A 36-7-47 29	3691B 36-7-47 30	3692A 36-7-47 33	3692B 36-7-47 34	3693A 36-7-47 37	3693B 36-7-47 38	3694A 36-7-47 41	3694B 36-7-47 42	3695A 36-7-47 45
				SAMPLE ID/#														
DATE				02/03/88	03/16/88	03/16/88	02/03/88	03/16/88	03/16/88	03/15/88	03/15/88	03/10/88	03/10/88	03/10/88	03/10/88	03/16/88	03/16/88	03/15/88
TIME				10:16	10:43	10:55	11:22	11:46	11:55	10:57	11:07	11:20	11:27	11:53	12:03	13:07	13:17	11:37
SAMPLE TYPE		71999	0	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO
SITE TYPE I		99759	0	CSDT	BORE	BORE	CSDT	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE
SAMPLE DEPTH	FT	99758	0	0	11	15	0	10	14	0	0	0	4	0	4	0	4	0
SAMPLING TECHNIQUE		72005	0	G	S	S	G	S	S	S	S	S	S	S	S	S	S	S
INSTALLATION CODE	SAMPLE	99720	0	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK
MOISTURE	%WET WT	70320	I	10.8	16.9	14.0	10.0	13.4	16.1	9.6	8.4	11.0	5.8	12.9	8.5	9.1	5.2	13.6
CADMIUM	UG/G- DRY	1028	R9	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921
CHROMIUM	UG/G-DRY	99584	R9	<7.16	14.0	9.92	17.2	9.39	12.0	12.3	12.4	<7.16	<7.16	<7.16	12.2	11.0	8.95	17.6
COPPER	UG/G- DRY	1043	R9	21.4	13.0	16.9	19.6	10.2	14.8	10.7	9.25	8.33	8.80	8.86	13.1	14.1	9.10	12.9
LEAD	UG/G-DRY	1052	R9	80.6	<16.8	<16.8	21.5	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	71.9	<16.8	<16.8
ZINC	UG/G-DRY	1093	R9	252	48.0	56.5	150	39.6	57.0	44.0	40.9	34.6	34.8	32.3	50.4	131	35.6	55.6
ARSENIC	UG/G- DRY	1003	T9	<4.70	<4.70	<4.70	6.69	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	15.2	<4.70	<4.70
MERCURY	UG/G-DRY	71921	V9	0.105	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050
ALDRIN	UG/G-DRY	98356	Q9	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94
ATRAZINE	UG/G-DRY	98655	Q9	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73
CHLORDANE	UG/G-DRY	98361	Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
P-CLPHENYLMETHY-SULFIDE	UG/G-DRY	98653	Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	8.1	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
P-CLPHENYLMETHY-SULFOXIDE	UG/G-DRY	98654	Q9	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	3.6	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35
P-CLPHENYLMETHY-SULFONE	UG/G-DRY	98703	Q9	<0.29	1.5	1.5	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
DBCP(NEMAGON)	UG/G-DRY	98652	Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
DICYCLOPENTADIENE	UG/G-DRY	98651	Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
DDE,PP'	UG/G-DRY	98363	Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
DDT,PP'	UG/G-DRY	98364	Q9	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37
DIELDRIN	UG/G-DRY	98365	Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	1.6	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25

PARAMETERS	UNITS	STORET #	METHOD	SAMPLE ID/# 3689GRAB 36-7-47 21	3689A 36-7-47 22	3689B 36-7-47 23	3690GRAB 36-7-47 25	3690A 36-7-47 26	3690B 36-7-47 27	3691A 36-7-47 29	3691B 36-7-47 30	3692A 36-7-47 33	3692B 36-7-47 34	3693A 36-7-47 37	3693B 36-7-47 38	3694A 36-7-47 41	3694B 36-7-47 42	3695A 36-7-47 45
DATE				02/03/88	03/16/88	03/16/88	02/03/88	03/16/88	03/16/88	03/15/88	03/15/88	03/10/88	03/10/88	03/10/88	03/10/88	03/16/88	03/16/88	03/15/88
TIME				10:16	10:43	10:55	11:22	11:46	11:55	10:57	11:07	11:20	11:27	11:53	12:03	13:07	13:17	11:37
DIMP	UG/G-DRY	98645	Q9	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
1,4 DITHIANE	UG/G-DRY	98650	Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
DMMP	UG/G-DRY	98657	Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
ENDRIN	UG/G-DRY.	98369	Q9	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70
HEXACHLOROCYCLOPENTADIENE	UG/G-DRY	98647	Q9	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1
ISODRIN	UG/G-DRY	98649	Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
MALATHION	UG/G-DRY	98648	Q9	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59
1,4 OXATHIANE	UG/G-DRY	98644	Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
ETY'PARATHION	UG/G-DRY	98658	Q9	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63
SUPONA	UG/G-DRY	98656	Q9	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49
VAPONA	UG/G-DRY	98646	Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
DICYCLOPENTADIENE	UG/G-DRY	98651	W9			<0.27			<0.27									
TRANS-1,2-DICHLOROETHENE	UG/G-DRY	98687	W9			<0.25			<0.25									
ETHYLBENZENE	UG/G-DRY	98688	W9			<0.25			<0.25									
METHYLENE CHLORIDE	UG/G-DRY	98689	W9			<0.25			0.69									
TETRACHLOROETHENE	UG/G-DRY	98690	W9			<0.25			<0.25									
TOLUENE	UG/G-DRY	98691	W9			<0.25			<0.25									
1,1,1-TRICHLOROETHANE	UG/G-DRY	98692	W9			<0.25			<0.25									
1,1,2-TRICHLOROETHANE	UG/G-DRY	98693	W9			<0.25			<0.25									
TRICHLOROETHENE	UG/G-DRY	98694	W9			<0.25			<0.25									
M-XYLENE	UG/G-DRY	98695	W9			<0.25			<0.25									
MIBK	UG/G-DRY	98696	W9			<0.50			<0.50									
DMDS	UG/G-DRY	98697	W9			<0.25			<0.25									
BENZENE	UG/G-DRY	98699	W9			<0.25			<0.25									

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3689GRAB 36-7-47 21	3689A 36-7-47 22	3689B 36-7-47 23	3690GRAB 36-7-47 25	3690A 36-7-47 26	3690B 36-7-47 27	3691A 36-7-47 29	3691B 36-7-47 30	3692A 36-7-47 33	3692B 36-7-47 34	3693A 36-7-47 37	3693B 36-7-47 38	3694A 36-7-47 41	3694B 36-7-47 42	3695A 36-7-47 45
DATE			02/03/88	03/16/88	03/16/88	02/03/88	03/16/88	03/16/88	03/15/88	03/15/88	03/10/88	03/10/88	03/10/88	03/10/88	03/16/88	03/16/88	03/15/88
TIME			10:16	10:43	10:55	11:22	11:46	11:55	10:57	11:07	11:20	11:27	11:53	12:03	13:07	13:17	11:37
UNK600		90600				2											
	UG/G	Q9															
UNK609		90609															
	UG/G	Q9															
UNK618		90618	0.8														
	UG/G	Q9															
UNK619		90619															
	UG/G	Q9															
UNK626		90626															
	UG/G	Q9															
UNK631		90631															
	UG/G	Q9															
UNK633		90633	0.9														
	UG/G	Q9															
UNK160		90160															
	UG/G	W9															
UNK615		90615						0.9									
	UG/G	Q9															
UNK543		90543															
	UG/G	Q9															
UNK559		90559															
	UG/G	Q9															
UNK573		90573															
	UG/G	Q9															
UNK576		90576				7											
	UG/G	Q9															
UNK578		90578				6											
	UG/G	Q9															
UNK622		90622															
	UG/G	Q9															
UNK624		90624	1.0														
	UG/G	Q9															
UNK628		90628	1														
	UG/G	Q9															
UNK632		90632	1	2													
	UG/G	Q9															
UNK634		90634															
	UG/G	Q9															
UNK636		90636	1		3												
	UG/G	Q9															
UNK640		90640	1														
	UG/G	Q9															
UNK644		90644															
	UG/G	Q9															
UNK649		90649															
	UG/G	Q9															
UNK655		90655															
	UG/G	Q9															

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3689GRAB 36-7-47 21	3689A 36-7-47 22	3689B 36-7-47 23	3690GRAB 36-7-47 25	3690A 36-7-47 26	3690B 36-7-47 27	3691A 36-7-47 29	3691B 36-7-47 30	3692A 36-7-47 33	3692B 36-7-47 34	3693A 36-7-47 37	3693B 36-7-47 38	3694A 36-7-47 41	3694B 36-7-47 42	3695A 36-7-47 45
DATE			02/03/88	03/16/88	03/16/88	02/03/88	03/16/88	03/16/88	03/15/88	03/15/88	03/10/88	03/10/88	03/10/88	03/10/88	03/16/88	03/16/88	03/15/88
TIME			10:16	10:43	10:55	11:22	11:46	11:55	10:57	11:07	11:20	11:27	11:53	12:03	13:07	13:17	11:37
UNK577		90577															
	UG/G	Q9															
UNK594		90594				2											
	UG/G	Q9															
UNK629		90629															
	UG/G	Q9															
UNK630		90630															
	UG/G	Q9															
UNK637		90637															
	UG/G	Q9															
UNK555		90555															
	UG/G	Q9															
UNK556		90556															
	UG/G	Q9															
UNK582		90582	4														
	UG/G	Q9															
UNK617		90617	2														
	UG/G	Q9															
UNK161		90161			3												
	UG/G	W9															
UNK541		90541			1.0												
	UG/G	Q9															
UNK551		90551				0.9											
	UG/G	Q9															
UNK591		90591				0.9											
	UG/G	Q9															
UNK595		90595				3											
	UG/G	Q9															
UNK518		90518													1		
	UG/G	Q9															
UNK661		90661													0.8		
	UG/G	Q9															
UNK523		90523															
	UG/G	Q9															
UNK616		90616															
	UG/G	Q9															
UNK673		90673															
	UG/G	Q9															
UNK550		90550															
	UG/G	Q9															
UNK572		90572															
	UG/G	Q9															
UNK627		90627															
	UG/G	Q9															
UNK526		90526							9								
	UG/G	Q9															
UNK614		90614	1		0.9												
	UG/G	Q9															

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3689GRAB 36-7-47 21	3689A 36-7-47 22	3689B 36-7-47 23	3690GRAB 36-7-47 25	3690A 36-7-47 26	3690B 36-7-47 27	3691A 36-7-47 29	3691B 36-7-47 30	3692A 36-7-47 33	3692B 36-7-47 34	3693A 36-7-47 37	3693B 36-7-47 38	3694A 36-7-47 41	3694B 36-7-47 42	3695A 36-7-47 45
DATE			02/03/88	03/16/88	03/16/88	02/03/88	03/16/88	03/16/88	03/15/88	03/15/88	03/10/88	03/10/88	03/10/88	03/10/88	03 16 88	03/16/88	03 15/88
TIME			10:16	10:43	10:55	11:22	11:46	11:55	10:57	11:07	11:20	11:27	11:53	12:03	13:07	13:17	11:37
UNK635	UG/G	90635 Q9															

PARAMETERS	STORET #	SAMPLE ID/# 3695B	3696A	3696B	3697A	3697B	3698A	3698B	3699A	3699B	3700A	3701A	3702A	3703A	3704A	3705A
UNITS	METHOD	36-7-47 46	36-7-47 49	36-7-47 50	36-7-47 53	36-7-47 54	36-7-47 57	36-7-47 58	36-7-47 61	36-7-47 62	36-7-47 65	36-7-47 67	36-7-47 69	36-7-47 71	36-7-47 73	36-7-47 75
DATE		03/15/88	03/14/88	03/14/88	03/15/88	03/15/88	03/16/88	03/16/88	03/15/88	03/15/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88
TIME		11:45	08:40	08:52	12:53	13:00	12:30	12:40	12:15	12:23	09:14	09:24	09:34	09:46	09:54	10:02
SAMPLE TYPE	71999 0	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO
SITE TYPE 1	99759 0	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE
SAMPLE DEPTH FT	99758 0	4	0	4	0	4	0	4	0	4	0	0	0	0	0	0
SAMPLING TECHNIQUE	72005 0	S	S	S	S	S	S	S	S	S	S	S	S	S	S	S
INSTALLATION CODE	99720 0	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK
SAMPLE MOISTURE %WET WT	70320 I	6.3	12.0	5.7	13.4	5.3	11.3	5.0	12.3	5.4	8.3	9.4	13.3	10.9	11.1	11.0
CADMIUM UG/G- DRY	1028 R9	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921
CHROMIUM UG/G-DRY	99584 R9	14.5	<7.16	11.9	9.27	<7.16	<7.16	9.50	<7.16	10.6	8.83	<7.16	11.1	11.8	15.5	10.3
COPPER UG/G- DRY	1043 R9	14.6	10.0	8.38	8.77	6.81	7.38	8.86	7.74	10.1	9.77	8.41	11.9	13.3	8.70	11.1
LEAD UG/G-DRY	1052 R9	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	22.5	<16.8	<16.8	<16.8
ZINC UG/G-DRY	1093 R9	56.1	39.0	36.3	45.1	32.5	33.5	39.2	35.3	43.8	36.9	28.3	48.9	45.9	33.4	51.6
ARSENIC UG/G- DRY	1003 T9	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70
MERCURY UG/G-DRY	71921 V9	2.08	<0.050	0.061	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	0.054	0.119	0.320	0.065	<0.050	<0.050
ALDRIN UG/G-DRY	98356 Q9	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94
ATRAZINE UG/G-DRY	98655 Q9	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73
CHLORDANE UG/G-DRY	98361 Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
P-CLPHENYLMETHY-SULFIDE UG/G-DRY	98653 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
P-CLPHENYLMETHY-SULFOXIDE UG/G DRY	98654 Q9	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35
P-CLPHENYLMETHY-SULFONE UG/G-DRY	98703 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
DBCP(NEMAGON) UG/G-DRY	98652 Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
DICYCLOPENTADIENE UG/G-DRY	98651 Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
DDE,PP' UG/G-DRY	98363 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
DDT,PP' UG/G-DRY	98364 Q9	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37
DIELDRIN UG/G-DRY	98365 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	4.4	<0.25	<0.25

PARAMETERS	UNITS	STORET #	METHOD	SAMPLE ID/# 3695B 36-7-47 46	3696A 36-7-47 49	3696B 36-7-47 50	3697A 36-7-47 53	3697B 36-7-47 54	3698A 36-7-47 57	3698B 36-7-47 58	3699A 36-7-47 61	3699B 36-7-47 62	3700A 36-7-47 65	3701A 36-7-47 67	3702A 36-7-47 69	3703A 36-7-47 71	3704A 36-7-47 73	3705A 36-7-47 75
DATE				03/15/88	03/14/88	03/14/88	03/15/88	03/15/88	03/16/88	03/16/88	03/15/88	03/15/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88
TIME				11:45	08:40	08:52	12:53	13:00	12:30	12:40	12:15	12:23	09:14	09:24	09:34	09:46	09:54	10:02
DIMP		98645		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
	UG/G-DRY		Q9															
1,4 DITHIANE		98650		<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
	UG/G-DRY		Q9															
DMMP		98657		<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
	UG/G-DRY		Q9															
ENDRIN		98369		<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70
	UG/G-DRY.		Q9															
HEXACHLOROCYCLOPENT-ADIENE	UG/G-DRY	98647	Q9	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1
ISODRIN		98649		<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
	UG/G-DRY		Q9															
MALATHION		98648		<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59
	UG/G-DRY		Q9															
1,4 OXATHIANE		98644		<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
	UG/G-DRY		Q9															
ETY'PARATHION		98658		<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63
	UG/G-DRY		Q9															
SUPONA		98656		<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49
	UG/G-DRY		Q9															
VAPONA		98646		<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
	UG/G-DRY		Q9															
DICYCLOPENTADIENE		98651																
	UG/G-DRY		W9															
TRANS-1,2-DICHLOROET HENE	UG/G-DRY	98687	W9															
ETHYLBENZENE		98688																
	UG/G-DRY		W9															
METHYLENE CHLORIDE		98689																
	UG/G-DRY		W9															
TETRACHLOROETHENE		98690																
	UG/G-DRY		W9															
TOLUENE		98691																
	UG/G-DRY		W9															
1,1,1-TRICHLORO-ETHANE	UG/G-DRY	98692	W9															
1,1,2-TRICHLORO-ETHANE	UG/G-DRY	98693	W9															
TRICHLOROETHENE		98694																
	UG/G-DRY		W9															
M-XYLENE		98695																
	UG/G-DRY		W9															
MIBK		98696																
	UG/G-DRY		W9															
DMDS		98697																
	UG/G-DRY		W9															
BENZENE		98699																
	UG/G-DRY		W9															

PARAMETERS	STORET #	SAMPLE ID/# 3695B 36-7-47 46	3696A 36-7-47 49	3696B 36-7-47 50	3697A 36-7-47 53	3697B 36-7-47 54	3698A 36-7-47 57	3698B 36-7-47 58	3699A 36-7-47 61	3699B 36-7-47 62	3700A 36-7-47 65	3701A 36-7-47 67	3702A 36-7-47 69	3703A 36-7-47 71	3704A 36-7-47 73	3705A 36-7-47 75
UNITS	METHOD															
DATE		03/15/88	03/14/88	03/14/88	03/15/88	03/15/88	03/16/88	03/16/88	03/15/88	03/15/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88
TIME		11:45	08:40	08:52	12:53	13:00	12:30	12:40	12:15	12:23	09:14	09:24	09:34	09:46	09:54	10:02
O-AND/OR P-XYLENE	98700															
UG/G-DRY	W9															
CARBON TETRACHLORIDE	98680															
UG/G-DRY	W9															
CHLOROBENZENE	98681															
UG/G-DRY	W9															
CHLOROFORM	98682															
UG/G-DRY	W9															
1,1-DICHLOROETHANE	98683															
UG/G-DRY	W9															
1,2-DICHLOROETHANE	98684															
UG/G-DRY	W9															
BICYCLOHEPTADIENE	98686															
UG/G-DRY	W9															
DBCP(NEMAGON)	98652															
UG/G-DRY	W9															
THIODIGLYCOL	99798															
UG/G	MM9															
CHLOROACETIC ACID	97285															
UG/G	MM9															
IMPA	97382															
UG/G	AAA9															
FLUOROACETIC ACID	97381															
UG/G	AAA9															
MPA	97383															
UG/G	AAA9															
UNK587	90587															
UG/G	Q9															
UNK588	90588															
UG/G	Q9															
UNK593	90593															
UG/G	Q9															
UNK602	90602															
UG/G	Q9															
UNK524	90524						9	8								
UG/G	Q9															
UNK549	90549															
UG/G	Q9															
UNK554	90554															
UG/G	Q9															
UNK558	90558						0.8									
UG/G	Q9															
UNK566	90566						1									
UG/G	Q9															
UNK586	90586						0.9									
UG/G	Q9															
UNK598	90598															
UG/G	Q9															

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3695B 36-7-47 46	3696A 36-7-47 49	3696B 36-7-47 50	3697A 36-7-47 53	3697B 36-7-47 54	3698A 36-7-47 57	3698B 36-7-47 58	3699A 36-7-47 61	3699B 36-7-47 62	3700A 36-7-47 65	3701A 36-7-47 67	3702A 36-7-47 69	3703A 36-7-47 71	3704A 36-7-47 73	3705A 36-7-47 75
DATE			03/15/88	03/14/88	03/14/88	03/15/88	03/15/88	03/16/88	03/16/88	03/15/88	03/15/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/86	03/09/86
TIME			11:45	08:40	08:52	12:53	13:00	12:30	12:40	12:15	12:23	09:14	09:24	09:34	09:46	09:54	10:02
UNK600		90600															
	UG/G	Q9															
UNK609		90609															
	UG/G	Q9															
UNK618		90618															
	UG/G	Q9															
UNK619		90619		0.9													
	UG/G	Q9															
UNK626		90626															
	UG/G	Q9															
UNK631		90631															
	UG/G	Q9															
UNK633		90633		0.8													
	UG/G	Q9															
UNK160		90160															
	UG/G	W9															
UNK615		90615															
	UG/G	Q9															
UNK543		90543															
	UG/G	Q9															
UNK559		90559															
	UG/G	Q9															
UNK573		90573															
	UG/G	Q9															
UNK576		90576															
	UG/G	Q9															
UNK578		90578															
	UG/G	Q9															
UNK622		90622															
	UG/G	Q9															
UNK624		90624															
	UG/G	Q9															
UNK628		90628		0.9													
	UG/G	Q9													2		
UNK632		90632															
	UG/G	Q9															
UNK634		90634															
	UG/G	Q9															
UNK636		90636										1				2	
	UG/G	Q9															
UNK640		90640															
	UG/G	Q9															
UNK644		90644															
	UG/G	Q9															
UNK649		90649															
	UG/G	Q9															
UNK655		90655															
	UG/G	Q9															

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PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3695B 36-7-47 46	3696A 36-7-47 49	3696B 36-7-47 50	3697A 36-7-47 53	3697B 36-7-47 54	3698A 36-7-47 57	3698B 36-7-47 58	3699A 36-7-47 61	3699B 36-7-47 62	3700A 36-7-47 65	3701A 36-7-47 67	3702A 36-7-47 69	3703A 36-7-47 71	3704A 36-7-47 73	3705A 36-7-47 75
DATE			03/15/88	03/14/88	03/14/88	03/15/88	03/15/88	03/16/88	03/16/88	03/15/88	03/15/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88
TIME			11:45	08:40	08:52	12:53	13:00	12:30	12:40	12:15	12:23	09:14	09:24	09:34	09:46	09:54	10:02
UNK577		90577															
	UG/G	Q9															
UNK594		90594															
	UG/G	Q9															
UNK629		90629															
	UG/G	Q9															
UNK630		90630															
	UG/G	Q9															
UNK637		90637															0.9
	UG/G	Q9															
UNK555		90555															
	UG/G	Q9															
UNK556		90556															
	UG/G	Q9															
UNK582		90582															
	UG/G	Q9															
UNK617		90617															
	UG/G	Q9															
UNK161		90161															
	UG/G	W9															
UNK541		90541															
	UG/G	Q9															
UNK551		90551															
	UG/G	Q9															
UNK591		90591															
	UG/G	Q9															
UNK595		90595															
	UG/G	Q9															
UNK518		90518															
	UG/G	Q9															
UNK661		90661															
	UG/G	Q9															
UNK523		90523													0.9		
	UG/G	Q9															
UNK616		90616													3		
	UG/G	Q9															
UNK673		90673													5		
	UG/G	Q9															
UNK550		90550															
	UG/G	Q9															
UNK572		90572															
	UG/G	Q9															
UNK627		90627															
	UG/G	Q9															
UNK526		90526		10		10											
	UG/G	Q9															
UNK614		90614															
	UG/G	Q9															

PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3695B 36-7-47 46	3696A 36-7-47 49	3696B 36-7-47 50	3697A 36-7-47 53	3697B 36-7-47 54	3698A 36-7-47 57	3698B 36-7-47 58	3699A 36-7-47 61	3699B 36-7-47 62	3700A 36-7-47 65	3701A 36-7-47 67	3702A 36-7-47 69	3703A 36-7-47 71	3704A 36-7-47 73	3705A 36-7-47 75
DATE			03/15/88	03/14/88	03/14/88	03/15/88	03/15/88	03/16/88	03/16/88	03/15/88	03/15/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88	03/09/88
TIME			11:45	08:40	08:52	12:53	13:00	12:30	12:40	12:15	12:23	09:14	09:24	09:34	09:46	09:54	10:02
UNK635	UG/G	90635 Q9				0.8	2										

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PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	STORET #	SAMPLE ID/# 3706A 36-7-47	3707A 36-7-47	3707B 36-7-47	3707C 36-7-47	3708A 36-7-47	3708B 36-7-47	3708C 36-7-47	3709A 36-7-47	3709B 36-7-47	3709C 36-7-47	3710A 36-7-47	3710B 36-7-47	3711A 36-7-47	3711B 36-7-47	3712A 36-7-47
UNITS	METHOD	77	79	80	81	84	85	86	89	90	91	94	95	97	98	100
DATE		03/09/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/14/88	03/14/88	03/14/88	03/14/88	03/14/88
TIME		10:10	08:21	08:33	08:46	09:20	09:29	09:46	10:20	10:30	10:47	13:18	13:22	12:49	12:54	13:43
SAMPLE TYPE	71999 0	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO
SITE TYPE 1	99759 0	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE	BORE
SAMPLE DEPTH FT	99758 0	0	0	4	9	0	4	9	0	4	9	0	2	0	2	0
SAMPLING TECHNIQUE	72005 0	S	S	S	S	S	S	S	S	S	S	S	S	S	S	S
INSTALLATION CODE SAMPLE	99720 0	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK
MOISTURE %WET WT	70320 1	11.7	8.7	8.6	8.1	10.6	6.7	9.5	12.2	8.2	12.2	8.3	3.0	10.4	5.8	11.5
CADMIUM UG/G- DRY	1028 R9	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921
CHROMIUM UG/G-DRY	99584 R9	12.1	<7.16	10.6	9.30	<7.16	9.08	16.0	11.5	14.9	11.3	<7.16	<7.16	11.0	10.3	<7.16
COPPER UG/G- DRY	1043 R9	9.26	6.97	9.35	10.7	11.2	9.76	12.3	14.3	13.7	13.1	8.77	6.54	9.06	7.80	7.52
LEAD UG/G-DRY	1052 R9	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8
ZINC UG/G-DRY	1093 R9	38.7	25.3	34.9	34.1	38.5	34.7	38.6	53.8	51.3	37.4	34.3	27.8	36.7	32.2	34.3
ARSENIC UG/G- DRY	1003 T9	6.15	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	<4.70	5.97					
MERCURY UG/G-DRY	71921 V9	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050	<0.050					
ALDRIN UG/G-DRY	98356 Q9	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94					
ATRAZINE UG/G-DRY	98655 Q9	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73					
CHLORDANE UG/G-DRY	98361 Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5					
P-CLPHENYLMETHY-SULFIDE UG/G-DRY	98653 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25					
P-CLPHENYLMETHY-SULFOXIDE UG/G-DRY	98654 Q9	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35					
P-CLPHENYLMETHY-SULFONE UG/G-DRY	98703 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29					
DBCP(NEMAGON) UG/G-DRY	98652 Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33					
DICYCLOPENTADIENE UG/G-DRY	98651 Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26					
DDE,PP' UG/G-DRY	98363 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29					
DDT,PP' UG/G-DRY	98364 Q9	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37					
DIELDRIN UG/G-DRY	98365 Q9	<0.25	<0.25	<0.25	<0.25	0.51	<0.25	<0.25	0.35	<0.25	<0.25					

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PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	UNITS	STORET #	METHOD	SAMPLE ID/# 3706A 36-7-47 77	3707A 36-7-47 79	3707B 36-7-47 80	3707C 36-7-47 81	3708A 36-7-47 84	3708B 36-7-47 85	3708C 36-7-47 86	3709A 36-7-47 89	3709B 36-7-47 90	3709C 36-7-47 91	3710A 36-7-47 94	3710B 36-7-47 95	3711A 36-7-47 97	3711B 36-7-47 98	3712A 36-7-47 100
DATE				03/09/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/14/88	03/14/88	03/14/88	03/14/88	03/14/88
TIME				10:10	08:21	08:33	08:46	09:20	09:29	09:46	10:20	10:30	10:47	13:18	13:22	12:49	12:54	13:43
DIMP		98645		<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50					
	UG/G-DRY		Q9															
1,4 DITHIANE		98650		<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25					
	UG/G-DRY		Q9															
DMMP		98657		<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5					
	UG/G-DRY		Q9															
ENDRIN		98369		<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70					
	UG/G-DRY.		Q9															
HEXACHLOROCYCLOPENT-		98647		<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1					
ADIENE	UG/G-DRY		Q9															
ISODRIN		98649		<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33					
	UG/G-DRY		Q9															
MALATHION		98648		<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59					
	UG/G-DRY		Q9															
1,4 OXATHIANE		98644		<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26					
	UG/G-DRY		Q9															
ETY'PARATHION		98658		<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63	<0.63					
	UG/G-DRY		Q9															
SUPONA		98656		<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49	<0.49					
	UG/G-DRY		Q9.															
VAPONA		98646		<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25					
	UG/G-DRY		Q9															
DICYCLOPENTADIENE		98651					<0.27			<0.27			<0.27					
	UG/G-DRY		W9															
TRANS-1,2-DICHLOROET		98687					<0.25			<0.25			<0.25					
HENE	UG/G-DRY		W9															
ETHYLBENZENE		98688					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
METHYLENE CHLORIDE		98689					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
TETRACHLOROETHENE		98690					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
TOLUENE		98691					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
1,1,1-TRICHLORO-		98692					<0.25			<0.25			<0.25					
ETHANE	UG/G-DRY		W9															
1,1,2-TRICHLORO-		98693					<0.25			<0.25			<0.25					
ETHANE	UG/G-DRY		W9															
TRICHLOROETHENE		98694					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
M-XYLENE		98695					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
MIBK		98696					<0.50			<0.50			<0.50					
	UG/G-DRY		W9															
DMDS		98697					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															
BENZENE		98699					<0.25			<0.25			<0.25					
	UG/G-DRY		W9															

PARAMETERS	STORET #	SAMPLE ID/# 3706A 36-7-47 77	3707A 36-7-47 79	3707B 36-7-47 80	3707C 36-7-47 81	3708A 36-7-47 84	3708B 36-7-47 85	3708C 36-7-47 86	3709A 36-7-47 89	3709B 36-7-47 90	3709C 36-7-47 91	3710A 36-7-47 94	3710B 36-7-47 95	3711A 36-7-47 97	3711B 36-7-47 98	3712A 36-7-47 100
UNITS	METHOD															
DATE		03/09/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/14/88	03/14/88	03/14/88	03/14/88	03/14/88
TIME		10:10	08:21	08:33	08:46	09:20	09:29	09:46	10:20	10:30	10:47	13:18	13:22	12:49	12:54	13:43
O-AND/OR P-XYLENE	98700				<0.50			<0.50			<0.50					
UG/G-DRY	W9															
CARBON TETRACHLORIDE	98680				<0.25			<0.25			<0.25					
UG/G-DRY	W9															
CHLOROBENZENE	98681				<0.25			<0.25			<0.25					
UG/G-DRY	W9															
CHLOROFORM	98682				<0.25			<0.25			<0.25					
UG/G-DRY	W9															
1,1-DICHLOROETHANE	98683				<0.25			<0.25			<0.25					
UG/G-DRY	W9															
1,2-DICHLOROETHANE	98684				<0.28			<0.28			<0.28					
UG/G-DRY	W9															
BICYCLOHEPTADIENE	98686				<0.25			<0.25			<0.25					
UG/G-DRY	W9															
DBCP(NEMAGON)	98652				<0.33			<0.33			<0.33					
UG/G-DRY	W9															
THIODIGLYCOL	99798															
UG/G	MM9															
CHLOROACETIC ACID	97285															
UG/G	MM9															
IMPA	97382															
UG/G	AAA9															
FLUOROACETIC ACID	97381															
UG/G	AAA9															
MPA	97383															
UG/G	AAA9															
UNK587	90587															
UG/G	Q9															
UNK588	90588															
UG/G	Q9															
UNK593	90593															
UG/G	Q9															
UNK602	90602															
UG/G	Q9															
UNK524	90524		IL													
UG/G	Q9															
UNK549	90549															
UG/G	Q9															
UNK554	90554															
UG/G	Q9															
UNK558	90558															
UG/G	Q9															
UNK566	90566															
UG/G	Q9															
UNK586	90586															
UG/G	Q9															
UNK598	90598															
UG/G	Q9															

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3706A 36-7-47 77	3707A 36-7-47 79	3707B 36-7-47 80	3707C 36-7-47 81	3708A 36-7-47 84	3708B 36-7-47 85	3708C 36-7-47 86	3709A 36-7-47 89	3709B 36-7-47 90	3709C 36-7-47 91	3710A 36-7-47 94	3710B 36-7-47 95	3711A 36-7-47 97	3711B 36-7-47 98	3712A 36-7-47 100
DATE			03/09/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/14/88	03/14/88	03/14/88	03/14/88	03/14/88
TIME			10:10	08:21	08:33	08:46	09:20	09:29	09:46	10:20	10:30	10:47	13:18	13:22	12:49	12:54	13:43
UNK600		90600															
	UG/G	Q9															
UNK609		90609															
	UG/G	Q9															
UNK618		90618															
	UG/G	Q9															
UNK619		90619															
	UG/G	Q9															
UNK626		90626				0.9											
	UG/G	Q9															
UNK631		90631															
	UG/G	Q9															
UNK633		90633															
	UG/G	Q9															
UNK160		90160				OK1.0			OK6								
	UG/G	W9															
UNK615		90615	1									0.9					
	UG/G	Q9															
UNK543		90543															
	UG/G	Q9															
UNK559		90559															
	UG/G	Q9															
UNK573		90573															
	UG/G	Q9															
UNK576		90576															
	UG/G	Q9															
UNK578		90578															
	UG/G	Q9															
UNK622		90622															
	UG/G	Q9															
UNK624		90624															
	UG/G	Q9															
UNK628		90628				0.8											
	UG/G	Q9															
UNK632		90632															
	UG/G	Q9															
UNK634		90634															
	UG/G	Q9															
UNK636		90636	1			4		0.8	1	8							
	UG/G	Q9															
UNK640		90640															
	UG/G	Q9															
UNK644		90644															
	UG/G	Q9															
UNK649		90649															
	UG/G	Q9															
UNK655		90655															
	UG/G	Q9															

PARAMETERS	UNITS	STORET # METHOD	SAMPLE ID/# 3706A 36-7-47 77	3707A 36-7-47 79	3707B 36-7-47 80	3707C 36-7-47 81	3708A 36-7-47 84	3708B 36-7-47 85	3708C 36-7-47 86	3709A 36-7-47 89	3709B 36-7-47 90	3709C 36-7-47 91	3710A 36-7-47 94	3710B 36-7-47 95	3711A 36-7-47 97	3711B 36-7-47 98	3712A 36-7-47 100
DATE			03/09/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/14/88	03/14/88	03/14/88	03/14/88	03/14/88
TIME			10:10	08:21	08:33	08:46	09:20	09:29	09:46	10:20	10:30	10:47	13:18	13:22	12:49	12:54	13:43
UNK577		90577															
	UG/G	Q9															
UNK594		90594															
	UG/G	Q9															
UNK629		90629															
	UG/G	Q9															
UNK630		90630															
	UG/G	Q9															
UNK637		90637															
	UG/G	Q9															
UNK555		90555															
	UG/G	Q9															
UNK556		90556															
	UG/G	Q9															
UNK582		90582															
	UG/G	Q9															
UNK617		90617															
	UG/G	Q9															
UNK161		90161															
	UG/G	W9															
UNK541		90541															
	UG/G	Q9															
UNK551		90551															
	UG/G	Q9															
UNK591		90591															
	UG/G	Q9															
UNK595		90595															
	UG/G	Q9															
UNK518		90518															
	UG/G	Q9															
UNK661		90661															
	UG/G	Q9															
UNK523		90523															
	UG/G	Q9															
UNK616		90616															
	UG/G	Q9															
UNK673		90673															
	UG/G	Q9															
UNK550		90550															
	UG/G	Q9															
UNK572		90572															
	UG/G	Q9															
UNK627		90627															
	UG/G	Q9															
UNK526		90526															
	UG/G	Q9															
UNK614		90614															
	UG/G	Q9															

PARAMETERS	UNITS	STORET # / METHOD	SAMPLE ID/# 3706A 36-7-47 77	3707A 36-7-47 79	3707B 36-7-47 80	3707C 36-7-47 81	3708A 36-7-47 84	3708B 36-7-47 85	3708C 36-7-47 86	3709A 36-7-47 89	3709B 36-7-47 90	3709C 36-7-47 91	3710A 36-7-47 94	3710B 36-7-47 95	3711A 36-7-47 97	3711B 36-7-47 98	3712A 36-7-47 100
DATE			03/09/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/10/88	03/14/88	03/14/88	03/14/88	03/14/88	03/14/88
TIME			10:10	08:21	08:33	08:46	09:20	09:29	09:46	10:20	10:30	10:47	13:18	13:22	12:49	12:54	13:43
UNK635		90635															
	UG/G	09															

SAMPLE ID/#

PARAMETERS	UNITS	STORET #	METHOD	3712B 36-7-47 101	3688A 36-7-47 103	3688B 36-7-47 104
DATE				03/14/88	03/16/88	03/16/88
TIME				13:46	09:40	09:51
SAMPLE TYPE		71999	0	SO	SO	SO
SITE TYPE 1		99759	0	BORE	BORE	BORE
SAMPLE DEPTH	FT	99758	0	2	9	13
SAMPLING TECHNIQUE		72005	0	S	S	S
INSTALLATION CODE	SAMPLE	99720	0	RK	RK	RK
MOISTURE	%WET WT	70320	1	5.6	15.6	21.6
CADMIUM	UG/G- DRY	1028	R9	<0.921	<0.921	<0.921
CHROMIUM	UG/G-DRY	99584	R9	11.3	<7.16	<7.16
COPPER	UG/G- DRY	1043	R9	9.00	11.4	40.5
LEAD	UG/G-DRY	1052	R9	<16.8	<16.8	<16.8
ZINC	UG/G-DRY	1093	R9	40.1	38.7	77.8
ARSENIC	UG/G- DRY	1003	T9		<4.70	<4.70
MERCURY	UG/G-DRY	71921	V9		0.125	<0.050
ALDRIN	UG/G-DRY	98356	Q9		<0.94	<0.94
ATRAZINE	UG/G-DRY	98655	Q9		<0.73	<0.73
CHLORDANE	UG/G-DRY	98361	Q9		<1.5	<1.5
P-CLPHENYLMETHY-SULFIDE	UG/G-DRY	98653	Q9		<0.25	<0.25
P-CLPHENYLMETHY-SULFOXIDE	UG/G-DRY	98654	Q9		<0.35	<0.35
P-CLPHENYLMETHY-SULFONE	UG/G-DRY	98703	Q9		<0.29	<0.29
DBCP(NEMAGON)	UG/G-DRY	98652	Q9		<0.33	<0.33
DICYCLOPENTADIENE	UG/G-DRY	98651	Q9		<0.26	<0.26
DDE,PP'	UG/G-DRY	98363	Q9		<0.29	<0.29
DDT,PP'	UG/G-DRY	98364	Q9		<0.37	<0.37
DIELDRIN	UG/G-DRY	98365	Q9		<0.25	<0.25

SAMPLE ID/#

PARAMETERS	UNITS	STORET # METHOD	3712B 36-7-47 101	3688A 36-7-47 103	3688B 36-7-47 104
DATE			03/14/88	03/16/88	03/16/88
TIME			13:46	09:40	09:51
DIMP		98645		<0.50	<0.50
	UG/G-DRY	Q9			
1,4 DITHIANE		98650		<0.25	<0.25
	UG/G-DRY	Q9			
DMMP		98657		<1.5	<1.5
	UG/G-DRY	Q9			
ENDRIN		98369		<0.70	<0.70
	UG/G-DRY.	Q9			
HEXACHLOROCYCLOPENT-		98647		<1.1	<1.1
ADIENE	UG/G-DRY	Q9			
ISODRIN		98649		<0.33	<0.33
	UG/G-DRY	Q9			
MALATHION		98648		<0.59	<0.59
	UG/G-DRY	Q9			
1,4 OXATHIANE		98644		<0.26	<0.26
	UG/G-DRY	Q9			
ETY'PARATHION		98658		<0.63	<0.63
	UG/G-DRY	Q9			
SUPONA		98656		<0.49	<0.49
	UG/G-DRY	Q9			
VAPONA		98646		<0.25	<0.25
	UG/G-DRY	Q9			
DICYCLOPENTADIENE		98651			<0.27
	UG/G-DRY	W9			
TRANS-1,2-DICHLOROET		98687			<0.25
HENE	UG/G-DRY	W9			
ETHYLBENZENE		98688			<0.25
	UG/G-DRY	W9			
METHYLENE CHLORIDE		98689			0.63
	UG/G-DRY	W9			
TETRACHLOROETHENE		98690			<0.25
	UG/G-DRY	W9			
TOLUENE		98691			<0.25
	UG/G-DRY	W9			
1,1,1-TRICHLORO-		98692			<0.25
ETHANE	UG/G-DRY	W9			
1,1,2-TRICHLORO-		98693			<0.25
ETHANE	UG/G-DRY	W9			
TRICHLOROETHENE		98694			<0.25
	UG/G-DRY	W9			
M-XYLENE		98695			<0.25
	UG/G-DRY	W9			
MIBK		98696			<0.50
	UG/G-DRY	W9			
DMDS		98697			<0.25
	UG/G-DRY	W9			
BENZENE		98699			<0.25
	UG/G-DRY	W9			

ENVIRONMENTAL SCIENCE & ENGINEERING 06/24/88
PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47/1
FIELD GROUP 36-7-47 PROJECT MANAGER BILL FRASER
ALL LAB COORDINATOR JOE VONDRICK

SAMPLE ID/#

PARAMETERS	UNITS	STORET # METHOD	3712B 36-7-47 101	3688A 36-7-47 103	3688B 36-7-47 104
DATE			03/14/88	03/16/88	03/16/88
TIME			13:46	09:40	09:51
UNK577		90577			
	UG/G	Q9			
UNK594		90594			
	UG/G	Q9			
UNK629		90629		20	
	UG/G	Q9			
UNK630		90630		200	
	UG/G	Q9			
UNK637		90637			
	UG/G	Q9			
UNK555		90555			
	UG/G	Q9			
UNK556		90556			
	UG/G	Q9			
UNK582		90582			
	UG/G	Q9			
UNK617		90617			
	UG/G	Q9			
UNK161		90161			
	UG/G	W9			
UNK541		90541			
	UG/G	Q9			
UNK551		90551			
	UG/G	Q9			
UNK591		90591			
	UG/G	Q9			
UNK595		90595			
	UG/G	Q9			
UNK518		90518			
	UG/G	Q9			
UNK661		90661			
	UG/G	Q9			
UNK523		90523			
	UG/G	Q9			
UNK616		90616			
	UG/G	Q9			
UNK673		90673			
	UG/G	Q9			
UNK550		90550		8	
	UG/G	Q9			
UNK572		90572		40	
	UG/G	Q9			
UNK627		90627		400	
	UG/G	Q9			
UNK526		90526			
	UG/G	Q9			
UNK614		90614			0.9
	UG/G	Q9			

SAMPLE ID/#

PARAMETERS	UNITS	STORET # METHOD	3712B 36-7-47 101	3688A 36-7-47 103	3688B 36-7-47 104
DATE			03/14/88	03/16/88	03/16/88
TIME			13:46	09:40	09:51
UNK635		90635			
	UG/G	Q9			

ENVIRONMENTAL SCIENCE & ENGINEERING 06/28/88 PAGE#1
PROJECT NUMBER 84936 0300 PROJECT NAME RMA TASK 1
FIELD GROUP TIMB PROJECT MANAGER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	STORET #	SAMPLE ID/# BLK TIMB 488	BLK T21MB2 2	BLK T47MB1 1	BLK T47MB1 6	BLK T47MB1 7	BLK T47MB1 8	BLK T21MB2 5	BLK T47MB1 21	BLK T47MB1 26	BLK T47MB1 28	BLK T47MB1 29	BLK TIMB 542	BLK T47MB1 41	BLK T47MB1 48	BLK T47MB1 49
UNITS	METHOD															
DATE		10/27/87	03/11/88	02/02/88	03/09/88	02/12/88	03/10/88	03/16/88	02/03/88	03/09/88	03/15/88	03/10/88	01/26/88	02/04/88	03/09/88	03/09/88
TIME																
SAMPLE TYPE	71999 0	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO	SO
SITE TYPE 1	99759 0	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB
SAMPLE DEPTH	99758	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
FT	0															
SAMPLING TECHNIQUE	72005 0	G	G	G	G	G	G	G	G	G	G	G	G	G	G	G
INSTALLATION CODE	99720	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK	RK
SAMPLE	0															
MOISTURE	70320	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4	2.4
%WET WT	0															
CADMIUM	1028	<0.921	<0.921	<0.921	<0.921	<0.921	<0.921									
UG/G- DRY	R9															
CHROMIUM	99584	12.1	14.3	10.4	9.99	<7.16	12.8									
UG/G-DRY	R9															
COPPER	1043	9.56	10.6	10.1	10.1	8.20	9.76									
UG/G- DRY	R9															
LEAD	1052	<16.8	<16.8	<16.8	<16.8	<16.8	<16.8									
UG/G-DRY	R9															
ZINC	1093	41.7	40.9	39.3	33.3	29.6	35.0									
UG/G-DRY	R9															
ARSENIC	1003							<4.70	5.59	<4.70	<4.70	<4.70				
UG/G- DRY	T9															
MERCURY	71921												<0.050	<0.050	<0.050	<0.050
UG/G-DRY	V9															
THIODIGLYCOL	99798															
UG/G	MM9															
CHLOROACETIC ACID	97285															
UG/G	MM9															

PARAMETERS	STORET #	SAMPLE ID/# BLK T47MB1	BLK T47MB1
UNITS	METHOD	50	61
DATE		03/16/88	02/02/88
TIME			
SAMPLE TYPE	71999	SO	SO
	0		
SITE TYPE 1	99759	QCMB	QCMB
	0		
SAMPLE DEPTH	99758	0	0
CM	0		
SAMPLING TECHNIQUE	72005	G	G
	0		
INSTALLATION CODE	99720	RK	RK
SAMPLE	0		
MOISTURE	70320	2.4	2.4
%WET WT	0		
CADMIUM	1028		
UG/G- DRY	R9		
CHROMIUM	99584		
UG/G-DRY	R9		
COPPER	1043		
UG/G- DRY	R9		
LEAD	1052		
UG/G-DRY	R9		
ZINC	1093		
UG/G-DRY	R9		
ARSENIC	1003		
UG/G- DRY	T9		
MERCURY	71921	<0.050	
UG/G-DRY	V9		
THIODIGLYCOL	99798		<2.55
UG/G	MM9		
CHLOROACETIC ACID	97285		<18.0
UG/G	MM9		

ENVIRONMENTAL SCIENCE & ENGINEERING 06/28/88 PAGE#1
PROJECT NUMBER 84936 0300 PROJECT NAME RMA TASK I
FIELD GROUP TIMB PROJECT MANAGER
ALL LAB COORDINATOR JOE VONDRICK

PARAMETERS	STORET #	BLK TIMB 411	BLK TIMB 412	BLK T47MB1 85	BLK T47MB1 86	BLK T47MB1 87	BLK T47MB1 88	SAMPLE ID/# BLK T47MB1 89
UNITS	METHOD							
DATE		02/02/88	02/02/88	03/09/88	03/10/88	03/09/88	03/14/88	03/16/88
TIME					09:15		10:20	
SAMPLE TYPE	71999 0	SO	SO	SO	SO	SO	SO	SO
SITE TYPE 1	99759 0	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB	QCMB
SAMPLE DEPTH	99758 0	0	0	0	0	0	0	0
FT								
SAMPLING TECHNIQUE	72005 0	G	G	G	G	G	G	G
INSTALLATION CODE	99720 0	RK	RK	RK	RK	RK	RK	RK
SAMPLE								
MOISTURE	70320 0	2.4	2.4	2.4	2.4	2.4	2.4	2.4
%WET WT								
ALDRIN	98356 Q9	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94	<0.94
UG/G-DRY								
ATRAZINE	98655 Q9	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73	<0.73
UG/G-DRY								
CHLORDANE	98361 Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
UG/G-DRY								
P-CLPHENYLMETHY-	98653 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
SULFIDE UG/G-DRY								
P-CLPHENYLMETHY-	98654 Q9	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35	<0.35
SULFOXIDE UG/G-DRY								
P-CLPHENYLMETHY-	98703 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
SULFONE UG/G-DRY								
DBCP(NEMAGON)	98652 Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
UG/G-DRY								
DICYCLOPENTADIENE	98651 Q9	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26	<0.26
UG/G-DRY								
DDE,PP'	98363 Q9	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29	<0.29
UG/G-DRY								
DDT,PP'	98364 Q9	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37	<0.37
UG/G-DRY								
DIELDRIN	98365 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
UG/G-DRY								
DIMP	98645 Q9	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50	<0.50
UG/G-DRY								
1,4 DITHIANE	98650 Q9	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25	<0.25
UG/G-DRY								
DMMP	98657 Q9	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5	<1.5
UG/G-DRY								
ENDRIN	98369 Q9	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70	<0.70
UG/G-DRY.								
HEXACHLOROCYCLOPENT-	98647 Q9	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1	<1.1
ADIENE UG/G-DRY								
ISODRIN	98649 Q9	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33	<0.33
UG/G-DRY								
MALATHION	98648 Q9	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59	<0.59
UG/G-DRY								

ENVIRONMENTAL SCIENCE & ENGINEERING 06/28/88 PAGE#1
PROJECT NUMBER 88425 0000 PROJECT NAME RMA TASK 47
FIELD GROUP T47MB1 PROJECT MANAGER
ALL LAB COORDINATOR JOE VONDRICK

SAMPLE ID/#

PARAMETERS	STORET #	BLK T47MB1	BLK T47MB1	BLK T47MB1	BLK T47MB1
UNITS	METHOD	227	228	229	230
DATE		03/10/88	03/14/88	03/15/88	03/16/88
TIME					
SAMPLE TYPE	71999 0	SO	SO	SO	SO
SITE TYPE 1	99759 0	QCMB	QCMB	QCMB	QCMB
SAMPLE DEPTH	99758 0	0	0	0	0
FT					
SAMPLING TECHNIQUE	72005 0	G	G	G	G
INSTALLATION CODE	99720 0	RK	RK	RK	RK
SAMPLE					
MOISTURE	70320 0	2.4	2.4	2.4	2.4
%WET WT					
DICYCLOPENTADIENE	98651 W9	<0.27	<0.27	<0.27	<0.27
UG/G-DRY					
TRANS-1,2-DICHLOROETHENE	98687 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
ETHYLBENZENE	98688 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
METHYLENE CHLORIDE	98689 W9	0.94	0.99	0.40	<0.25
UG/G-DRY					
TETRACHLOROETHENE	98690 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
TOLUENE	98691 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
1,1,1-TRICHLOROETHANE	98692 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
1,1,2-TRICHLOROETHANE	98693 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
TRICHLOROETHENE	98694 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
M-XYLENE	98695 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
MIBK	98696 W9	<0.50	<0.50	<0.50	<0.50
UG/G-DRY					
DMDS	98697 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
BENZENE	98699 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
O-AND/OR P-XYLENE	98700 W9	<0.50	<0.50	<0.50	<0.50
UG/G-DRY					
CARBON TETRACHLORIDE	98680 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
CHLOROBENZENE	98681 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
CHLOROFORM	98682 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					
1,1-DICHLOROETHANE	98683 W9	<0.25	<0.25	<0.25	<0.25
UG/G-DRY					

SAMPLE ID/#

PARAMETERS	STORET #	BLK T47MB1	BLK T47MB1	BLK T47MB1	BLK T47MB1
UNITS	METHOD	227	228	229	230
DATE		03/10/88	03/14/88	03/15/88	03/16/88
TIME					
1,2-DICHLOROETHANE	98684	<0.28	<0.28	<0.28	<0.28
UG/G-DRY	W9				
BICYCLOHEPTADIENE	98686	<0.25	<0.25	<0.25	<0.25
UG/G-DRY	W9				
DBCP(NEMAGON)	98652	<0.33	<0.33	<0.33	<0.33
UG/G-DRY	W9				
UNK160	90160	10			
UG/G	W9				
UNK159	90159		5	100	5
UG/G	W9				